Submitted to:

Massachusetts Department of Environmental Protection Southeast Regional Office

January 12, 2011

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Massachusetts Department of Environmental Protection Southeast Regional Office Bureau of Waste Site Cleanup 20 Riverside Drive Lakeville, MA 02347

Re: Class A-2 Response Action Outcome Statement and Phase V Completion Statement
Walpole Park South
Walpole, Massachusetts
RTN 4-3021915

Dear Sir/Madam:

In accordance with the requirements of 310 CMR 40.000, and on behalf of Walpole Park South Trust, Tetra Tech, Inc. d/b/a Tetra Tech Rizzo is submitting this Class A-2 Response Action Outcome (RAO) Statement and Phase V Completion Statement for the Disposal Site identified by Release Tracking Number (RTN) 4-3021915, Walpole Park South (the Site). A Method 2 Human Health and Environmental Risk Characterization was conducted to evaluate the risk posed by the contaminants of concern at the Site. The results of the risk characterization concluded that a condition of No Significant Risk has been achieved for the Disposal Site, and an Activity and Use Limitation (AUL) is not required.

This report has been prepared in accordance with the Massachusetts Contingency Plan (MCP) 310 CMR 40.0000 and is subject to the Limitations and Conditions presented in Appendix A. The original DEP transmittal forms for this report (BWSC-104 and BWSC-108) were submitted electronically via eDEP.

Please contact us if you have any questions or comments.

Very truly yours,

Raymond C. Johnson, P.G., L.S.P. Senior Vice President

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1.0 Introduction

Tetra Tech, Inc. d/b/a Tetra Tech Rizzo (TTR) has prepared this Class A-2 Response Action Outcome (RAO) Statement and Phase V Completion Statement on behalf of Walpole Park South Trust for the Disposal Site identified by the Massachusetts Department of Environmental Protection under Release Tracking Number (RTN) 4-3021915 (the Site). The Site is known as Walpole Park South and is generally located at the northwest quadrant of the intersection of Providence Highway (Route 1) and Pine Street in Walpole, Massachusetts.

Multiple phases of subsurface investigations have been implemented at the Site, as described in the Phase II – Comprehensive Site Assessment (July 2006), Phase III – Remedial Action Plan (July 2006), Phase IV – Remedy Implementation Plan (July 2007), Phase IV Completion Report and Remedy Operation Status Submittal (July 2009) and Phase V – Remedy Operation Status Reports #1 and #2 submitted in February and August 2010, respectively. The Phase II report concluded that a condition of No Significant Risk did not exist in 2006 because of the presence of bromodichloromethane and chloroform in MW-2 and RIZ-3, chloroform in MW-3 and lead in MW-9. The installation of additional groundwater monitoring wells, and collection of six (6) rounds of groundwater samples since submission of the Phase II report generated additional groundwater data, indicative of current and recent conditions at the Site, which was used for preparation of a Method 2 risk characterization. Based on the results of the testing done since submission of the Phase II report and the conclusions of the risk characterization presented herein, a condition of No Significant Risk has been achieved at the Site and an Activity and Use Limitation (AUL) is not required to restrict or limit human exposures.

This report presents a description of the releases, a summary of historic and recent assessment of soil and groundwater conditions at the Site, a representativeness evaluation and data usability assessment, and a summary of the risk characterization.

This Class A-2 RAO Statement/Phase V Completion Statement has been prepared in accordance with the Massachusetts Contingency Plan (MCP), 310 CMR 40.0000, and is subject to the LSP Statement of Limitations and Conditions included in Appendix A. Copies of the DEP transmittal forms (BWSC-104 and BWSC-108) are included in Appendix F.

2.0 Disposal Site Information

In accordance with the MCP (310 CMR 40.1056) the following sections provide the general Disposal Site information relevant to this Class A-2 RAO and Phase V Completion Statement.

2.1 Disposal Site Description

The Site encompasses approximately 54 acres of land, located at the intersection of US Route 1 and Pine Street in Walpole, Massachusetts, as shown on Figure 1. The Site is divided into eight lots, seven of which contain buildings occupied by office and warehouse space that are leased to commercial and/or light industrial businesses. The current configuration of the Site and the

configuration of the individual building lots are depicted on Figure 2. An access road, Walpole Park South Drive, crosses the Site from Route 1 along the northeast boundary of Walpole Park South, to Pine Street on the southern boundary of the property. The buildings, driveways and parking areas cover the majority of the Site. The remainder of the property consists of landscaped areas adjacent to the buildings, wooded land and unpaved open areas. Prior to construction of the existing buildings the Site was vacant land, portions of which were reportedly used as a gravel pit. Development of the Site and building construction commenced in 1986.

The property is abutted to the north by vacant wooded land, to the west by single-family residences, to the south by Pine Street, across which are commercial properties including a truck repair and painting facility, and to the east by Route 1, across which are commercial and industrial properties. A Site Locus map is included as Figure 1, the Disposal Site boundaries and Site features are depicted on Figure 2 and the MassGIS Map is included as Figure 3.

The nearest surface water features to the Site are the Goldwater Farm Pond and School Meadow Brook. Goldwater Farm Pond is located approximately 650 feet south of the Site, and School Meadow Brook flows in a north and northwesterly direction downstream from the pond. The Site and surrounding area receive water from the Town of Walpole municipal water system, and are also serviced by public sanitary sewer service. No private drinking water wells are known or anticipated to exist within 500 ft of the Site. There is no known Protected Open Space within 500 ft of the Site. There are no known institutions (facilities with overnight housing such as a hospital, health care facility, orphanage, nursing home, convalescent home, educational facility, or correctional institution) located within 500 ft of the Site. Human receptors within 500 feet of the Disposal Site include workers at the Site and abutting properties, and residents of single family residences located west of the Site.

2.2 Release Description and Regulatory History

2.2.1 Background and Site History

Prior to construction of the existing buildings the Site was vacant land, portions of which were used as a gravel pit. Development of the Site and building construction commenced in 1986. Buildings have been constructed on the Site lots on an intermittent basis since 1986, and currently one lot remains undeveloped.

In compliance with requirements issued by the Walpole Board of Health (BOH), seven groundwater monitoring wells were installed at the Site in December 1986 by Carr Research Laboratory (Carr), and annual groundwater monitoring was performed on a limited number of the installed wells. In the late 1990's it was discovered that the reported results from the annual BOH sampling were occasionally exceeding the then current Massachusetts DEP reportable concentrations for GW-1 areas (RCGW-1 standards). Two RTNs were issued based on reported concentrations of compounds identified in groundwater at the Site.

RTN 3-19859 was issued in 2000 based on sampling performed in April 1999 and April 2000 when elevated concentrations of chloroform and bromodichloromethane were reported in

groundwater samples collected from monitoring wells along the western boundary of the Site. A Class A-2 RAO Statement was submitted by Carr Research Laboratory on July 25, 2001, indicating that a Permanent Solution had been achieved for this RTN. The RAO indicated that the source of the bromodichloromethane was treatment chemicals (brominating tablets) used in a residential swimming pool located on an upgradient property. The source of the chloroform was identified as a reaction between chlorine used for swimming pool disinfection at the same residence, and septage from the septic tank and leach field at the residence.

RTN 3-21915 was issued in April 2002 when it was noted that the lead concentrations reported by the laboratory for groundwater samples collected from monitoring wells MW-3 and MW-6 were 59 micrograms per liter (μ g/l) and 23 μ g/l, respectively; concentrations which exceeded the then applicable MCP reportable concentration for groundwater category RCGW-1 (20 μ g/l). To further evaluate this condition confirmatory groundwater sampling was performed in May 2002. The results of this sampling indicated lead concentrations in samples collected from MW-3 and MW-6 of 46 μ g/l and 18 μ g/l, respectively. Based on these sampling results it was concluded that the detected lead concentrations represented a 120-day notification condition under the MCP. Therefore, a RNF was submitted to the DEP on July 2, 2002. In response to the notification, DEP issued a Notice of Responsibility (NOR) on August 15, 2002 and assigned RTN 3-21915 to the reported release. Between September 2000 and February 2006 additional groundwater monitoring wells were installed on several occasions to replace damaged wells and/or to provide sufficient coverage to assess groundwater conditions at the Site.

Based on historic monitoring results and the investigations implemented after submission of the RNF for RTN 3-21915, a Phase I – Initial Site Investigation (Phase I) report and Tier Classification was prepared by GHC and submitted to DEP in June 2004. The Phase I report concluded that the nature and extent of contamination does not exhibit a regular pattern, relative to both the locations of wells in which levels exceed MCP Method 1 GW-1 standards and the detection of compounds over time. Based on the Numerical Ranking Scoresheet (NRS) prepared by GHC, the Site was classified as a Tier IB Disposal Site. An evaluation performed as part of the Phase I investigation concluded that the identified Site conditions did not represent an Imminent Hazard, indicating that implementation of an Immediate Response Action was not necessary. In an internal memorandum dated July 9, 2004, the DEP Drinking Water Program (DWP) found that "the groundwater contamination levels at the site are all low, compared to most waste sites," and "heavy metals have fairly low mobility in groundwater." DEP concluded "the site does not appear to pose a threat to the Walpole municipal wells, because of the low groundwater contamination levels and the distance from the site to the wells."

2.2.2 Phase II – Comprehensive Site Assessment

Following the Phase I investigation Rizzo Associates, Inc. (now Tetra Tech Rizzo) was retained to perform the Phase II – Comprehensive Site Assessment. The purpose of the Phase II investigation was to obtain data to characterize the nature and extent of releases of oil and/or hazardous materials (OHM) at the Site, quantify the risks posed by such releases, and assess the need to conduct further remedial actions at the Site.

The Phase II Investigation included the installation of 7 soil borings and completion of 3 of the soil borings as groundwater monitoring wells; sampling and analysis of soil and groundwater from select soil borings and monitoring wells; surveying to determine the locations and relative elevations of each newly installed monitoring well casing; gauging of groundwater elevations to evaluate the groundwater flow direction and prepare a potentiometric surface map; and hydraulic conductivity testing of three groundwater monitoring wells.

Laboratory analysis of seven (7) soil samples collected during the Phase II investigation did not detect VOCs or total metals at concentrations exceeding the applicable MCP Method 1 standards. Laboratory analysis of 45 groundwater samples collected over four Phase II sampling rounds identified the presence of volatile organic compounds (VOCs) and/or dissolved metals at concentrations greater than one or more of the applicable MCP Method 1 standards in four (4) wells (MW-2, MW-3, MW-9 and RIZ-3), all located near the perimeter of the Site and downgradient from potential off-site sources of groundwater contamination. The compounds exceeding the MCP Method 1 standards included bromodichloromethane, chloroform and lead; however, no on-site source was been identified for these compounds and the dissolved metals concentrations were inconsistent over the four sampling events. Lead was identified at a concentration greater than the method detection limit in well MW-9 in only one of the four groundwater sampling events. Because the locations and concentrations of the identified compounds were not consistent throughout the four groundwater sampling events there is not a clearly definable plume for the identified dissolved metals compounds.

The presence of bromodichloromethane and/or chloroform along the eastern boundary of the Site, specifically in monitoring wells MW-2, MW-3 and RIZ-3, suggested impacts from releases of chlorinated or brominated water. Historically identified concentrations of bromodichloromethane and chloroform along the western boundary of the Site, associated with RTN 3-19859, were attributed to a release of swimming pool water from the property upgradient from well MW-6. The chlorinated and/or brominated water was expected to have been reacting with the naturally occurring organic material in the sandy soils of the Site to form bromodichloromethane and chloroform. A similar reaction was also likely taking place along the eastern boundary of the Site. Potential sources for the chlorinated or brominated water are leaks in municipal water pipes, fire hydrant flushing, or infiltrating rainwater mixed with roadway deicing chemicals such as calcium chloride or sodium chloride. Based on the decreasing concentrations of bromodichloromethane and chloroform moving downgradient across the Site from well MW-2 to well RIZ-3 to well MW-3, the source of chlorinated or brominated water was likely upgradient from well MW-2, in or on the southeastern side of Route 1.

Since the concentrations of several compounds in groundwater exceeded the MCP Method 1 GW-1 and GW-3 standards, and the Site is located within a Zone II for a public water supply, the risk characterization performed during the Phase II concluded that a condition of No Significant Risk to human health and the environment had not been achieved at the Site for groundwater at that time.

2.2.3 Phase III – Remedial Action Plan

The Phase III evaluation included a review of alternative methods for treatment of groundwater to evaluate whether there were one or more financially and technically feasible remedial alternatives that could be implemented to reduce risk at the Site to a level where a Permanent Solution could be achieved, and the selection of an alternative for implementation. For the evaluation of remedial alternatives feasible technologies were considered based on their ability to address the conditions identified on the Site.

An initial screening evaluated nine remedial alternatives based on their ability to target these contaminant characteristics and the subsurface conditions at the Site. Alternatives evaluated during the initial screening included groundwater pump-and-treat, in-situ chemical oxidation, permeable reactive barriers, bioremediation/bio-barrier, electrical resistance heating, surfactant flushing, air sparging and vapor extraction, soil excavation and disposal/treatment, and MNA. Of these alternatives, groundwater pump-and-treat and MNA were selected for detailed evaluation.

The detailed evaluation compared the two remedial alternatives noted above in greater detail, based on the following criteria: effectiveness, reliability, difficulty, costs, risks, benefits and time for implementation. Based on the detailed evaluation, MNA was selected as the remedy for the Site.

At the time the Phase III was submitted it was anticipated that while the Phase IV was not due until July 26, 2007, MNA monitoring would commence in September or October 2006. However, because of unwillingness on the part of the Town of Walpole to cooperate with the responsible party relative to the installation of additional monitoring wells needed to implement the MNA, installation of the additional monitoring wells and commencement of the MNA sampling program was delayed until December 2007.

2.2.4 Phase IV - Remedy Implementation Plan

Based on the results of the Phase III, the Phase IV report indicated that MNA would be implemented at the Site to further evaluate groundwater conditions over time. The Phase IV report also noted that although MNA was identified as the appropriate remedial action for the Site, it may be determined that implementation of one or more other technologies should be considered as additional data on groundwater conditions was developed. In that case, feasible remedial alternatives would be evaluated and a determination made of whether the approach should be modified or changed. If changes to the remedial program were determined to be applicable, supplemental Phase III and Phase IV reports would be prepared to discuss the selection (Phase III) and design (Phase IV) of the remedy or remedies. The selection of MNA was appropriate given the sporadic and intermittent detection of metals or VOCs at concentrations exceeding applicable MCP standards, the absence of an identifiable source(s) of the detected compounds, and the likely off-site origin of those compounds. The proposed design for MNA included the installation of additional monitoring wells upgradient from the Site, and the collection of groundwater samples from the new wells and selected existing on-site monitoring wells.

2.2.5 Phase IV Completion Statement

Installation of the three additional monitoring wells, near the upgradient (southern) property line for Walpole Park South, was completed in December 2007. Wells were installed at two locations on Walpole Park South property, adjacent to Pine Street, and at one location on MHD property within the "jug handle" intersection of Route 1 southbound and Pine Street. The drilling locations were accessed using an all-terrain vehicle mounted hollow stem auger drilling rig, equipped with the capability to drill into bedrock, since the locations on Walpole Park South property were not accessible to conventional truck-mounted drilling equipment.

Four rounds of groundwater samples were collected during Phase IV. Based on the results of those samplings and previous sampling results for the Site, it was concluded that the requirements for a Class A or Class B RAO had not been and that ongoing monitoring would be performed to further characterize groundwater conditions over time. Specifically, it was indicated in the Phase IV Completion Statement that groundwater samples would be collected at approximately six month intervals for analysis for VOCs and metals. Details of the well installation and groundwater sampling were detailed in the Phase IV Completion Statement dated July 28, 2009.

2.2.6 Phase V - Remedy Operation Status

Phase V groundwater sampling was performed at the Site in December 2009 and June 2010. The results of those sampling events were discussed in Phase V – Remedy Operation Status Reports #1 and #2, submitted in February and August 2010, respectively. The results of four rounds of Phase IV groundwater monitoring and two rounds of Phase V groundwater monitoring did not detect VOCs or metals at concentrations exceeding the applicable MCP RCGW-1 reportable concentrations or the MCP Method 1 GW-1, GW-2 (where applicable) or GW-3 standards. The results of six (6) rounds of groundwater sampling performed over the period from December 2007 through June 2010 are summarized on Table 3.

3.0 Nature and Extent of Contamination

Based on the extensive subsurface testing completed to date, it does not appear that there is a specific on-site source of the identified compounds, historic data does not indicate a clearly definable plume, and recent results have shown the few detected analytes at concentrations well below the lower of the MCP GW-1 or GW-3 standards. Therefore, since contamination is not currently present on the Site, characterization of the nature and extent of contamination is not possible or required. However, historic groundwater monitoring results, including Phase II data and monitoring since submission of the Phase II report, are discussed in the following section to provide a general context for purposes of this RAO Statement/Phase V Completion Statement.

3.1 Extent of Groundwater Contamination

3.1.1 Phase II Groundwater Analysis Results

For the 45 groundwater samples that were submitted for laboratory analysis over four sampling rounds as a part of the Phase II investigation, VOCs and or dissolved metals concentrations greater than one or more of the applicable MCP Method 1 standards were reported in 4 wells. Compounds exceeding the MCP Method 1 standards included bromodichloromethane, chloroform and lead; however, no on-site source was identified for these compounds. Reported dissolved metals concentrations were inconsistent over the sampling events, and lead was identified at a concentration greater than the method detection limit in well MW-9 in only one of the four groundwater sampling events. Because the locations and concentrations of the identified compounds were not consistent throughout the four groundwater sampling events there is not a clearly definable dissolved metals plume.

The presence of bromodichloromethane and/or chloroform along the eastern boundary of the Site, specifically in monitoring wells MW-2, MW-3 and RIZ-3, suggested impacts from releases of chlorinated or brominated water. Historically identified concentrations of bromodichloromethane and chloroform along the western boundary of the Site, associated with RTN 3-19859, were attributed to a release of swimming pool water from the property upgradient from well MW-6. The chlorinated and/or brominated water was expected to have been reacting with the naturally occurring organic material in the sandy soils of the Site to form bromodichloromethane and chloroform. A similar reaction was also likely taking place along the eastern boundary of the Site. Potential sources for the chlorinated or brominated water are leaks in municipal water pipes, fire hydrant flushing, or infiltrating rainwater mixed with roadway deicing chemicals such as calcium chloride or sodium chloride. Based on the decreasing concentrations of bromodichloromethane and chloroform moving downgradient across the Site from well MW-2 to well RIZ-3 to well MW-3, the source of chlorinated or brominated water was likely upgradient from well MW-2, in or on the southeastern side of Route 1. We note that the Water Quality Reports for 2004 and 2005 issued by the Walpole Sewer & Water Department indicated the presence of bromodichloromethane and chloroform in samples collected from the municipal water system, and indicate that these compounds are a "by-product of drinking water disinfection." The generation of these compounds from chlorination of groundwater is a typical occurrence in this area of Massachusetts. Based on the data from monitoring wells MW-2, MW-3, RIZ-3 and RIZ-9, the Phase II report concluded that a condition of No Significant Risk did not exist in 2006.

3.1.2 Phase IV and Phase V Groundwater Analysis Results

Since submission of the Phase II report six (6) rounds of groundwater samples have been collected from monitoring wells MW-3, MW-9, RIZ-3, RIZ-8, RIZ-9 and RIZ-10. During that same period four (4) rounds of samples were collected from MW-2; samples could not be collected from that well during the December 2007 and December 2009 sampling events because the well was covered with snow piles and could not be located. Two rounds of groundwater samples were also collected from monitoring well RIZ-8S; samples were not collected at a

greater frequency from that well because during the other sampling events the well was dry or there was an insufficient saturated thickness to permit collection of water samples. These wells were selected to further evaluate conditions characterized during the Phase II investigations, to evaluate conditions near the upgradient property boundary, and to provide general coverage of the Site. Laboratory analysis of the samples collected during Phase IV and Phase V activities at the Site indicated low to non-detectable concentrations of the compounds which were the basis for the conclusions of the risk characterization in the Phase II report, and demonstrate that the historic elevated analysis results are likely related to off-site impacts that are intermittent and sporadic.

3.1.3 Summary

Historically, metals including antimony, arsenic, cadmium, chromium, and lead were identified at elevated concentrations at the Site. However, no metals have been reported at concentrations greater than the current Method 1 GW-1 standards since April 2006. In addition, when antimony was detected it was reported in multiple wells during a single groundwater sampling event but not reported at elevated concentrations in subsequent samplings of the same wells. The infrequent detections of antimony over the extensive sampling duration at the Site suggests that sampling and or laboratory error may account for the reported results, they may be naturally occurring impacts that vary in concentration over time, or related to intermittent off-site releases that migrate through the Site.

Lead is the only dissolved metal that has been identified in the groundwater at the Site on a somewhat regular basis; however, even lead concentrations have not been identified consistently enough to create plume maps or suggest an on-site source. Since April 1991, based on a combination of historical data and the groundwater sampling performed as a part of the Phase II investigation and post-Phase II monitoring, lead has been reported at concentrations greater than the current Method 1 GW-1 standard (15 μ g/L) three times in well MW-3 (22 to 59 μ g/L), twice in MW-9 (23 to 35 μ g/L) and once each in MW-2 (18 μ g/L), MW-6 (18 μ g/L) and MW-8 (26 μ g/L). Lead has not been reported above the lower of the MCP Method 1 GW-1 or GW-3 standard in the six sampling events performed since December 2007. Of these wells only well MW-3 is located on the downgradient side of the Site, indicating that an upgradient source or sources may be a significant contributor to the elevated lead concentrations on the Site.

None of the reported groundwater concentrations were detected at levels exceeding their respective Upper Concentration Limits (UCLs). Groundwater analytical data supporting this Class B-2 RAO is presented in Table 3. Laboratory certificates of analysis for groundwater samples collected as part of the MCP investigations since submission of the Phase II report are presented in Appendix E.

4.0 Representativeness Evaluation and Data Usability Assessment

A Representativeness Evaluation and Data Usability Assessment (REDUA) was conducted in support of this RAO Statement in accordance with 310 CMR 40.1056(2)(k) and DEP Policy

#WSC-07-350, "MCP Representativeness Evaluations and Data Usability Assessments" dated September 19, 2007. The Representativeness Evaluation documents the adequacy of the spatial and temporal data sets used to support the RAO. The Data Usability Assessment documents that the data relied upon are scientifically valid and defensible, and of sufficient accuracy, precision and completeness to support the RAO.

4.1 Representativeness Evaluation

The Representativeness Evaluation demonstrates the adequacy of the cumulative data set to sufficiently characterize conditions at the Disposal Site and supports the Conceptual Site Model. The evaluation includes a description of the Conceptual Site Model, use of field screening data; sampling rationale; number, spatial distribution and sampling procedures; temporal distribution of samples; data gaps; inconsistency and uncertainty and representativeness information. The components of the Representativeness Evaluation in support of this RAO are discussed in the sections below.

4.1.1 Conceptual Site Model

The Disposal Site is currently occupied by an industrial park with multiple buildings occupied by office, warehouse and distribution operations, paved parking lots, roadways and open space/landscaped areas. Laboratory analysis results of historic groundwater samples identified low concentrations of metals and VOCs. The monitoring has not identified a source or definable plume. The locations at which positive analytical results have been reported, and the intermittent and sporadic nature of the results, demonstrate that the historic elevated analysis results are likely related to off-site sources.

The topography of the Site slopes from southwest to northeast, from upland areas on the south and southwest side of Pine Street to wetlands and School Meadow Brook located north and northeast of the Site. The depth to bedrock ranges from 12 feet to greater than 50 feet below the ground surface, and the unconsolidated deposits overlying bedrock consist primarily of fine to coarse sand with gravel and boulders, with interbedded finer layers observed in some borings. Groundwater is present in overburden throughout most of the Site, although in some areas along the south and southwest property boundary, where shallow bedrock was observed, the overburden saturated thickness is limited and some monitoring wells are observed to periodically be dry. Groundwater at the site originates from migration from off-site upgradient locations, and from infiltration of precipitation at the Site, and flows from southwest to northeast consistent with topography and the groundwater discharge areas to the north and northeast (School Meadow Brook and associated wetlands).

4.1.2 Use of Field/Screening Data

The selection of laboratory analytical methods for the samples collected during and since the Phase II field investigations, which commenced in 2005, was based on the results of earlier sampling which indicated that VOCs and metals were the compounds of concern at the Site.

Field screening techniques, including PID headspace screening and field observations of soil characteristics and odors, were used in conjunction with existing analytical data to assist with the selection of soil samples for laboratory analysis.

4.1.3 Sampling Locations and Depths

Because of the absence of an identified release or source area, and the sporadic and intermittent presence of target compounds, sampling was conducted at locations throughout the Site to provide a representative characterization of subsurface conditions.

Our review indicates that the sampling locations and depths are sufficient to delineate Disposal Site boundaries, identify background COC concentrations, calculate EPCs, identify Hot Spots, identify exposure pathways and receptors, and assess human health and environmental risk at the Site.

4.1.4 Sampling Density, Spatial Distribution, Collection Methods and Handling

Soil and groundwater samples used to support this RAO were collected based on the layout and use of the Site and to provide general site-wide coverage. Soil samples included in the data set that supports this Class A-2 RAO Statement/Phase V Completion Statement were collected as discrete samples of subsurface materials. Groundwater samples were collected from monitoring wells screened across and below the water table using modified US Environmental Protection Agency (EPA) low flow sampling protocol which provides a representative sample of the static groundwater. Soil and groundwater samples collected to support this RAO were placed on ice and were transported to a Massachusetts certified laboratory under chains-of-custody.

The spatial distribution of the samples used for the calculation of EPCs used in the Risk Characterization is shown on Figure 2. A total of 30 soil samples, collected from 19 soil boring locations throughout the Site, were used to calculate the soil EPCs. Soil sample depths were selected based upon Field/Screening data, visual and olfactory observations and to provide general coverage of the fill materials encountered. A total of 48 groundwater samples collected from 9 monitoring wells during the period between December 2007 and June 2009 were used to calculate groundwater EPCs.

It is our opinion that the sampling density and spatial distribution of samples collected is consistent with the Conceptual Site Model and are representative of current Disposal Site conditions. We conclude that the sample collection and handling methods, sampling density and spatial distribution of soil and groundwater samples site-wide are sufficient to support the Class A-2 RAO Statement/Phase V Completion Statement.

4.1.5 Temporal Distribution of Samples

Soil and groundwater samples collected to support this RAO were collected at various times during the period from May 2003 to December 2007 for soil and December 2007 to June 2009

for groundwater. Therefore, we conclude that the temporal distribution of soil and groundwater sample collection is sufficient to demonstrate that no ongoing or uncontrolled source of contamination remains, that concentrations are stable and below the lower of the MCP GW-1 or GW-3 standards, and that EPCs accurately reflect Disposal Site concentrations.

4.1.6 Completeness

An extensive data set of 48 groundwater and 30 soil analytical data points has been used in support of this RAO. No data gaps were identified. Our review has concluded that the data set is sufficiently complete to support the RAO due to the extensive data set, sample density and distribution, and consistency with the Conceptual Site Model.

4.1.7 Inconsistency and Uncertainty

Based on the number of samples, historic site use, absence of on-site sources, and the consistency of the data set over multiple sampling events, no inconsistencies or uncertainty were identified in the data set.

4.1.8 Representativeness of the Data Set

Laboratory certificates of analysis for the groundwater samples collected since submission of the Phase II report are included in Appendix E and summarized on Table 3.

4.2 Data Usability Assessment

The Data Usability Assessment consists of an analytical and field component. The field component evaluates the sampling collection procedure to ensure a sample is representative of the sampling point upon delivery to the laboratory. The analytical component evaluates whether analytical data points are scientifically valid and defensible, and that a sufficient level of precision, accuracy, and sensitivity has been achieved. According to DEP Policy, the rigorousness of the Data Usability Assessment should be "proportional to the complexity of the project and the ramifications of risk-related decisions associated with the interpretation of the data."

As part of an effort to enhance the quality and consistency of analytical data supporting MCP submittals, the DEP has published the Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data (WSC-CAM-VII A). The original version of this document was released on May 21, 2004 and a revised version became effective on July 1, 2010. The purpose of the document is to provide quality assurance/quality control (QA/QC) guidance regarding the acquisition and reporting of analytical data, including regulatory performance standards and agency expectations for MCP data submittal. To facilitate the application of the performance standards, the DEP has published a Compendium of Analytical Methods (CAM) which provides a series of recommended protocols for the acquisition, analyses, and reporting of

analytical data in support of MCP decisions. Use of these procedures provides an LSP with "Presumptive Certainty" of data acceptance by the DEP.

Validation of the laboratory data included a review of field and laboratory quality control samples as applicable, including:

- CAM Compliance status and Presumptive Certainty of data;
- field sample custody, preservation, and analytical holding times;
- trip blanks to assess whether there may be false positive contamination during sample acquisition and/or storage;
- evaluation of surrogate recoveries to assess laboratory matrix effects and data accuracy;
- laboratory method blanks, surrogate spike recoveries (organics only), laboratory control sample (LCS) recoveries, MS/MSD recoveries, and laboratory sample duplicates and relative percent difference (RPD);
- laboratory quantitation (detection) limits; and
- miscellaneous observations.

A table summarizing the evaluation of analytical data points as part of the Data Usability Assessment is provided in Appendix B. The table lists the sample name, sample parameters, date sampled, sample matrix, CAM Compliance status and notes any qualifications or exceptions affecting data quality. The Data Usability Assessment provides justification as to why such analytical data are considered acceptable to support the RAO.

4.2.1 Analytical Data Usability Assessment

Soil data used in support of this RAO Statement were collected from May 2003 through December 2007, and groundwater data was collected from December 2007 to June 2009. The majority of the data is considered CAM Compliant, meaning that analytical results were (1) determined using an MCP Analytical Method detailed in CAM; (2) comply with method-specific QC requirements specified in CAM; (3) are reported with narration of method-specific performance standard deficiencies; and (4) are reported with required deliverables specified in CAM. The only data that is "non-CAM" is the June 2010 groundwater samples analyzed for VOCs by EPA Method 524. Specifically, the CAM method for VOCs is EPA 8260; however, Method 524 was used for groundwater VOC analysis to be consistent with the analytical method specified in the requirements for the annual groundwater monitoring performed at the Site in accordance with an agreement between the property owner and the Walpole Board of Health, and because Method 524 offers lower detection limits than Method 8260. CAM Compliant data have "Presumptive Certainty" and are of known accuracy, precision and sensitivity. Data that does not have "Presumptive Certainty" status may still be relied upon for purposes of the

conclusions of the RAO Statement provided further evaluation of the data concludes that it is of suitable quality to meet the requirements of the MCP.

A discussion of analytical data issues as identified by the Laboratory Case Narrative is included in Appendix B. In general, CAM Compliant data were found to have adequate sensitivity in detection levels compared to project-specific objectives. In some instances, surrogate recoveries, relative percent differences, matrix spike or method blank samples suggested that results may be biased low or detected outside of quality control limits. Most of these situations involved COCs that were consistently non-detect in samples collected throughout the Site or COCs that were materially contributing to the level of risk at the Site. The "non-CAM" data discussed above was also found to have adequate sensitivity in detection levels compared to project specific objectives, and this data is of suitable quality for use in the Risk Characterization and as a basis for this RAO Statement/Phase V Completion Statement.

In conclusion, limitations, uncertainties and qualifications associated with analytical deficiencies do not affect the overall accuracy, precision or sensitivity of the analytical data to the extent that the validity of the data would be jeopardized. No data were rejected based on gross failure criteria.

4.2.2 Field Data Usability Assessment

Appropriate sampling and handling methods were employed in the field during the collection of sample media. Suitable sampling containers were used, samples were properly preserved, stored within acceptable temperature ranges, and hold times were met for all samples. Analytical data were found to be consistent with field observations and screening data.

The field component of the data usability assessment indicates that sample collection, handling and analytical procedures were followed in a manner that is consistent with QC protocols required by the MCP. Based on our overall evaluation of the data, field sampling and handling methods are not anticipated to impact the field accuracy and precision of the data to the extent that the validity of the data would be jeopardized.

4.3 Conclusions

A Representativeness Evaluation and Data Usability Assessment was conducted in accordance with 310 CMR 40.1056(2) (k). It is our opinion that the data set relied upon to support this RAO is representative with regard to the spatial and temporal distribution of sampling points; is scientifically valid and defensible; and is of sufficient accuracy, precision and completeness to support the RAO Statement.

5.0 Method 2 Human Health and Environmental Risk Characterization

The following section presents a Method 2 Risk Characterization to evaluate the potential risks posed to human health, public welfare, public safety and the environment by concentrations of compounds of concern (COCs) that have been detected in Site soils and groundwater. This Risk Characterization was conducted in conformance with the requirements of the Massachusetts Contingency Plan (MCP), 310 CMR 40.0000. The DEP guidance document for risk characterization, *Guidance for Disposal Site Risk Characterization In Support of the Massachusetts Contingency Plan* (July 1995), has been followed in this analysis.

5.1 Method Selection

The MCP defines three methods for risk characterization: Methods 1, 2, and 3. We selected Method 2 as the appropriate method for characterization of risk because Method 1 standards have not been published by the DEP for isoproplybenzene and p-isopropyltoluene in soil. Method 2 allows for a relatively comprehensive, rapid evaluation of risk by comparison of Exposure Point Concentrations (EPCs) to standards published by DEP, and development of supplemental standards for compounds that do not have them or modification of standards with Site-specific information. Method 1 and 2 standards incorporate conservative assumptions for both contaminant transport and exposure, resulting in an overall conservative analysis. For the Site COCs that do not have Method 1 standards, we generated Method 2 standards for following the risk characterization guidance.

5.2 Soil and Groundwater Characterization

According to the criteria outlined in 310 CMR 40.0361, Site soils are classified as S-1 if the site is located within 500 feet of residentially zoned land, a school, playground, recreational area or park. Portions of the Site are located within 500 feet of residential property and, therefore, those areas are classified as category S-1.

The Massachusetts Geographic Information System (MassGIS) map of the Site area shows that the Site is located within a DEP Approved Zone II area and an EPA Sole Source Aquifer. According to the DEP, this means the Site is located within the "area of an aquifer which contributes water to a well under the most severe pumping and recharge conditions that can be realistically anticipated, as approved by the Department's Division of Water supply pursuant to 310 CMR 22.00." Thus, groundwater at the Site is classified as GW-1. Groundwater at the Site is also classified as GW-2 since there are occupied buildings on the Site and the depth to the water table is less than 15 feet. Additionally, groundwater at the Site is classified as GW-3, since, by definition, all groundwater in the Commonwealth of Massachusetts is considered GW-3. Therefore, we compare groundwater analytical results to GW-1, GW-2 and GW-3 standards.

5.3 Exposure Point Concentrations

EPCs represent the estimated concentrations of compounds of concern (COCs) to which a receptor may be exposed at the point of exposure. In keeping with DEP guidance, this characterization assumes that contaminant concentrations on the Site remain unchanged. Thus, we do not consider any mitigating factors resulting over the course of time (such as biodegradation). The risk characterization also assumes that for compounds detected at least once above detection limit, samples reported as not detected (ND) by the laboratory are assumed to have a concentration of one-half of the method detection limit (MDL) for that sample.

Data used in this risk characterization include analytical results for soil samples collected from the Site from May 2003 to December 2007. We also used groundwater data collected at the Site in December 2007, April, May, November and December 2008, June and December 2009, and June 2010. Tables 2 and 3 present the soil and groundwater data used in the risk characterization, respectively. As shown in Table 3, thallium was detected above the laboratory method detection limit for the sample collected in November 2008 from monitoring well RIZ-10, located on MassDOT property upgradient from the Site, at a concentration of 11.6 µg/l. Since thallium had not been recently detected at the Site, it was believed that the positive result for thallium could be related to sampling or analytical error. Therefore, the well was re-sampled in December 2008 to evaluate whether the original thallium result was representative of conditions in the area of RIZ-10. Analysis of the second sample did not detect thallium, confirming that the original result was not representative of groundwater conditions in the area of RIZ-10. Based on this we consider the positive thallium result for the sample collected from RIZ-10 in November 2008 to be anomalous and therefore, thallium is not considered a COC at the Site and is not carried through the risk characterization. In this risk characterization, we evaluate risk from all other compounds that were detected in the Site soil and groundwater.

For groundwater exposures, data from each monitoring well were evaluated as separate exposure points as required by the guidance for a Method 2 risk characterization under the MCP. This approach provides a conservative, health-protective assessment of risk. For soil exposures, we estimated four separate sets of soil EPCs (EPC-1 through EPC-4) for the Site based on the depth at which the soil samples were collected. EPC-1 is estimated as the average concentrations of COCs detected in soil samples collected 0 to 3 feet below the ground surface (bgs). EPC-2 is estimated as the average concentrations of COCs in soil samples collected within 3 to 7 feet bgs. EPC-3 is calculated as the average concentrations of COCs in soil samples collected from 8 to 15bgs. EPC-4 is estimated as the average concentrations of COCs in soil samples collected from 15 to 50 feet bgs.

The soil and groundwater EPCs estimated for the Site are presented in Tables 2 and 3, respectively.

6.0 Risk Characterization

Under the MCP Method 2 risk characterization a condition of "No Significant Risk" (NSR) of harm to human health, public welfare and the environment shall exist if each of the EPCs are

equal to or less than their applicable Method 2 standards, and there are no risks to public safety. This section presents a comparison of risk conditions with reference standards.

We note that Tables 2 and 3 indicate Method 2 standards for many compounds. However, as noted previously, only two compounds detected in soil do not have MCP Method 1 standards and a Method 2 risk characterization was performed to develop comparison standards for those compounds. In accordance with DEP Risk Assessment Guidance all comparison standards presented in the Tables must be identified as Method 2 standards, but for all but two compounds the listed comparison standards are the MCP Method 1 standards.

6.1.1 Risk of Harm to Health, Public Welfare and the Environment

As noted above, two soil COCs do not have Method 1 standards (isopropylbenzene, p-isopropyltoluene). These COCs are identified as tentatively identified compounds (TICs). The mass of these compounds is included in the mass of the aliphatic and aromatic hydrocarbons analyzed as extractable petroleum hydrocarbons (EPH). Therefore, the estimated Site soil EPCs of these compounds are compared to EPH (C_9 - C_{18} aliphatics) standards.

In Table 2, soil EPCs are compared with Method 2 S-1/GW-1, S-1/GW-2 and S-1/GW-3 standards. As shown in the table, the estimated soil EPCs are below the established Method 2 standards. In Table 3, groundwater EPCs are compared to Method 2 GW-1, GW-2 and GW-3 standards. The groundwater EPCs are below the applicable Method 2 standards. Therefore, we conclude that a condition of NSR to human health, the environment and public welfare exists at the Site for the conditions evaluated.

6.1.2 Risk of Harm to Public Safety

Threats to public safety include physical conditions and chemical agents that may cause bodily harm or injury (e.g. burns or fractures) as opposed to illness. There are no open pits, lagoons, drums, dangerous structures, or other apparent threats to public safety and no danger of fire or explosion from the conditions evaluated in this report. Thus, we find a condition of NSR of harm to public safety exists for the conditions observed at the Site.

6.2 Risk Characterization Conclusions

Based on the results of this Method 2 Risk Characterization, a condition of NSR of harm to human health, the environment, public welfare and public safety exists at the Site for the conditions evaluated. Since a condition of NSR exists without restriction of Site uses or activities the Site qualifies for a Class A-2 Response Action Outcome Statement.

7.0 Feasibility of Achieving Background

As part of this Response Action Outcome Statement the feasibility of reducing concentrations of OHM to achieve or approach background was evaluated in accordance with 310 CMR 40.0860

and the DEP Guidance Document "Conducting Feasibility Evaluations Under the MCP" (Policy #WSC-04-160) dated July 16, 2004. The evaluation criteria applicable to background feasibility include whether the benefits justify the costs or risks associated with the actions taken to achieve background; whether the technology needed to achieve or approach background exists; whether there are individuals with the necessary expertise to conduct the necessary remedial actions; or, if the selected alternative is off-site disposal, whether permitted facilities exist to accept the contaminated media. In accordance with DEP guidance, a finding of infeasibility based on any one of the criteria above is sufficient to conclude that achieving or approaching background is infeasible.

Extensive monitoring at the Site has not identified an on-site source of the compounds detected in groundwater or a definable contaminant plume. Further, the intermittent and sporadic nature of positive analytical results observed over many years indicates that COCs detected in soil and groundwater are either naturally occurring or related to off-site impacts that are migrating through the Site. Naturally-occurring compounds are by definition background that do not require further evaluation, and compounds related to off-site impacts migrating through the Site are beyond the control of the property owner. Therefore, achieving background is not feasible for the conditions identified at the Site.

8.0 Relationship of RAO to Other Disposal Site RAO(s)

310 CMR 40.1056(1) (d) requires a discussion of the relationship between this RAO and any other RAOs that have been filed for the Disposal Site. As discussed previously in this report, RTN 3-19859 was issued in 2000 based on sampling performed in April 1999 and April 2000 which detected elevated concentrations of chloroform and bromodichloromethane in groundwater samples collected near the western property boundary. A Class A-2 RAO Statement was submitted by Carr Research Laboratory on July 25, 2001, indicating that a Permanent Solution had been achieved for this RTN. The RAO indicated that the source of the bromodichloromethane was treatment chemicals (brominating tablets) used in a residential swimming pool located on an upgradient property. The source of the chloroform was identified as a reaction between chlorine used for swimming pool disinfection at the same residence and septage from the septic tank and leach field at the residence.

9.0 Phase V Completion Statement

As discussed herein, and documented in Phase V – ROS Status Reports submitted in February and August 2010, groundwater monitoring performed since submission of the Phase II report has demonstrated that groundwater conditions are at or approaching background, a condition of No Significant Risk has been achieved, an AUL is not needed to limit exposures or Site uses, and further monitoring is not required. Therefore, ROS is being terminated and conditions consistent with completion of the Phase V status of the Site have been achieved. The DEP Transmittal Form for Termination of ROS and Phase V Completion is in Appendix F.

10.0 Public Notifications

As stipulated by 310 CMR 40.1403(3) (f), and 40.1403(7), the following public involvement activities have been completed for the Site:

- Written notification to the parties identified on the Public Involvement Plan (PIP) mailing
 list was made at least 14 days in advance of the January 13, 2011 PIP meeting and an
 advertisement was published in the December 30, 2011 edition of The Walpole Times to
 notify the general public of the PIP meeting. Therefore, the advance notification
 requirements for a PIP were met.
- Letters have been sent notifying the Chief Municipal Officer and the Board of Health of the availability of this RAO Statement/Phase V Completion Statement.

Copies of these public involvement filings are included in Appendix C.

11.0 L.S.P. Opinion and Response Action Outcome Statement

Six rounds of groundwater sampling have been performed at the Site during the period from December 2007 to June 2009 to further evaluate groundwater conditions at monitoring wells where the Phase II report concluded a condition of Significant Risk existed n the July 26, 2006 Phase II report. Remedial actions at the Site were limited to Monitored Natural Attenuation as discussed in the Phase IV report. The results of the risk characterization concluded that a condition of "No Significant Risk" of harm to human health, the environment, public safety or public welfare exists for the current and anticipated future Site conditions and uses evaluated, and that an Activity and Use Limitation is not required to limit uses or exposures at the Site.

In accordance with the MCP, the category of an RAO achieved for a Disposal Site is established based upon the following factors: whether the site poses "No Significant Risk"; whether all Substantial Hazards posed by the Disposal Site have been eliminated; whether remedial actions were undertaken to achieve a level of "No Significant Risk"; whether one or more AULs are required to maintain a level of "No Significant Risk"; whether concentrations of oil or hazardous materials (OHM) at the site exceed the Upper Concentration Limits (UCLs) in soil and/or groundwater; and whether remedial actions have achieved background in accordance with the MCP. A comparison of these criteria to the conditions evaluated for the Disposal Site indicates that the requirements for a Class A-2 RAO have been met since a Permanent Solution has been achieved, OHM concentrations have not been reduced to background, an AUL is not required, and OHM do not exceed applicable UCLs [310 CMR 40.1036(3)]. The L.S.P. Statement of Limitations and Conditions, and copies of the DEP Transmittal Forms for this Class A-2 RAO Statement/Phase V Completion Statement, are included in Appendix A and Appendix F, respectively.

MCP. A comparison of these criteria to the conditions evaluated for the Disposal Site indicates that the requirements for a Class A-2 RAO have been met, since a Permanent Solution has been achieved; OHM concentrations have not been reduced to background; an AUL is not required and OHM do not exceed applicable UCLs [310 CMR 40.1036(3)]. The L.S.P. Statement of Limitations and Conditions, and a copy of the DEP Transmittal Forms for this Class A-2 RAO Statement/Phase V Completion Statement, are included in Appendix A and Appendix F, respectively.



Table 1. Groundwater Elevation Data - April 2010

| Monitoring Well I.D. | Relative Well Casing Elevation (ft) | Depth to Groundwater (ft) | Relative Groundwater Elevation (ft) |
|----------------------|-------------------------------------|---------------------------|-------------------------------------|
| MW-1 | 259.36 | 13.25 | 246.11 |
| MW-2 | 240.90 | 4.49 | 236.41 |
| MW-3 | 236.67 | 31.49 | 205.18 |
| MW-4 | 229.74 | 32.35 | 197.39 |
| MW-5S | 238.03 | 13.90 | 224.13 |
| MW-5D | 236.36 | 11.87 | 224.49 |
| MW-8 | 258.61 | 11.80 | 246.81 |
| MW-9 | 256.08 | 22.72 | 233.36 |
| GHC-1 | 241.95 | 4.73 | 237.22 |
| GHC-2 | 258.51 | 12.37 | 246.14 |
| GHC-3 | 252.40 | 13.51 | 238.89 |
| GHC-5 | 236.94 | 32.26 | 204.68 |
| GHC-6 | 236.01 | 1.35 | 234.66 |
| RIZ-1 | 239.60 | NM ⁶ | NM ⁶ |
| RIZ-2 | 234.94 | 4.85 | 230.09 |
| RIZ-3 | 241.52 | 9.10 | 232.42 |
| RIZ-8 | 265.52 | 20.17 | 245.35 |
| RIZ-8S | 265.38 | 19.61 | 245.77 |
| RIZ-9 | 246.69 | 5.85 | 240.84 |
| RIZ-10 | No survey data | 30.70 | No survey data |

Notes

- 1. Rim elevations for MW and GHC series wells from the "Groundwater Sampling Report, Winter-Spring 2004" Report, August 10, 2004
- 2. Rim elevations for RIZ-1 through RIZ-3 from April 10, 2006 Rizzo Associates survey
- 3. Rim elevations for RIZ-8 and RIZ-8S surveyed relative to MW-1, MW-8 and GHC-2 on April 25, 2008.
- 4. Rim elevations for RIZ-9 surveyed relative to RIZ-1 on April 25, 2008 $\,$
- 5. Depth to water measured during groundwater screening on April 6, 2010
- 6. RIZ-1 not gauged due to well being under water and water upwelling proximate to well.

| ocation: | Walpole Pk S. | | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S |
|-----------------------------|---------------|-------------|---------------|---------------|---------------|---------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| ample Name: | MW-3 (5ft) | MW-3 (50ft) | MW-1-S1-5 | MW-1-S2-10 | MW-1-S3-15 | MW-1-S4-20 | GHC-1 SS-1 | GHC-1 SS-3 | GHC-2 SS-1 | GHC-2 SS-4 | GHC-3 SS-1 | GHC-3 SS-3 | GHC-4 SS-1 | GHC-4 SS-3 | GHC-5 SS-1 | GHC-5 SS-4 | GHC-6 SS-1 | GHC-6 SS-2 | GHC-7 SS-1 | GHC-7 SS-4 |
| ample Depth: | 5' | 50' | 3'-5' | 8'-10' | 13'-15' | 18'-20' | 0-2' | 10'-12' | 0-2' | 15'-17' | 0-2' | 10'-12' | 0-2' | 10'-12' | 0-2' | 15'-17' | 0-2' | 5'-7' | 0-2' | 15'-17' |
| aboratory: | | | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum | Spectrum |
| aboratory I.D.: | | | AD93588 | AD93589 | AD93590 | AD93591 | SA07466-01 | SA07466-03 | SA07466-04 | SA07466-06 | SA07466-08 | SA07466-10 | SA07466-12 | SA07466-14 | SA07466-15 | SA07466-18 | SA07466-24 | SA07466-25 | SA07466-27 | SA07466-3 |
| ample Date: | 21-May-03 | 21-May-03 | 16-Jun-03 | 16-Jun-03 | 16-Jun-03 | 16-Jun-03 | 19-Jan-04 |
| onsultant: | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC | GHC |
| | | | | | | | 0.770 | <0.783 | <0.478 | 0.040 | 0.000 | 0.400 | 0.405 | 0.000 | 0.540 | 0.700 | 0.000 | 0.500 | <0.887 | 0.040 |
| etone nvlbenzene | | | | | | | <0.773 <0.0386 | <0.783 <0.0392 | <0.478 <0.0239 | <0.610 <0.0305 | <0.668 <0.0334 | <0.408 <0.0204 | <0.435 <0.0218 | <0.690 <0.0345 | <0.543 <0.0272 | <0.703 <0.0352 | <0.800 <0.0400 | <0.508 <0.0254 | <0.887 <0.0443 | <0.819 <0.0409 |
| opropylbenzene | | | | | | | < 0.0386 | <0.0392 | <0.0239 | <0.0305 | <0.0334 | <0.0204 | <0.0218 | <0.0345 | <0.0272 | <0.0352 | <0.0400 | <0.0254 | <0.0443 | <0.0409 |
| opropyltoluene, p- | | | | | | | <0.0386 | < 0.0392 | <0.0239 | <0.0305 | <0.0334 | <0.0204 | <0.0218 | <0.0345 | <0.0272 | <0.0352 | <0.0400 | <0.0254 | <0.0443 | <0.0409 |
| thylene chloride | | | | | | | | | | | | | | | | | | | | |
| chloromethane) | | | | | | | < 0.386 | < 0.392 | < 0.239 | < 0.305 | < 0.334 | < 0.204 | <0.218 | < 0.345 | < 0.272 | < 0.352 | < 0.400 | < 0.254 | < 0.443 | < 0.409 |
| oluene | | | | | | | <0.0386 | <0.0392 | <0.0239 | <0.0305 | < 0.0334 | <0.0204 | <0.0218 | < 0.0345 | <0.0272 | <0.0352 | <0.0400 | <0.0254 | <0.0443 | <0.0409 |
| senic. Total | | | | | | | 3.9 | 4.1 | <2.97 | <2.89 | <3.08 | <2.91 | <3.03 | <3.47 | <3.25 | <2.94 | <3.41 | <3.47 | <3.22 | <3.37 |
| arium, Total | | | | | | | 17.5 | 32.7 | 18.3 | 36.9 | 9.7 | 11.1 | 15.8 | 42.7 | 14.1 | 11.3 | 17.2 | 34.9 | 19.8 | 24.6 |
| ryllium, Total | | | | | | | 7.0 | 0.0 | 44.7 | 00.7 | 0.5 | 0.4 | 40.0 | 4 7 | 0.7 | 44.0 | 40.0 | - 0 | 0.4 | 40.7 |
| romium, Total ad. Total | 5.2 | 5.4 | 14.5 | 11.8 | 90.1 | 22.7 | 7.8 8.6 | 3.9 4.4 | 11.7 <1.98 | 30.7 18.8 | 6.5 3.1 | 8.1 <1.45 | 10.9 | 4.7 <4.62 | 9.7 <2.06 | 11.0 <1.47 | 10.0 4.2 | 5.6 5.0 | 9.4 8.7 | 12.7 <1.96 |
| ckel. Total | 5.2 | 5.4 | 14.5 | 11.0 | 90.1 | 22.1 | 0.0 | 4.4 | <1.90 | 10.0 | 3.1 | <1.45 | <1.61 | <4.02 | <2.06 | <1.47 | 4.2 | 5.0 | 0.7 | <1.90 |
| anadium | | | | | | | | | | | | | | | | | | | | |
| nc. Total | | | | | | | | | | | | | | | | | | | | |
| ercury, Total | | | | | | | <0.208 | <0.169 | <0.166 | <0.192 | <0.191 | <0.173 | <0.186 | <0.204 | <0.174 | <0.175 | <0.176 | <0.184 | <0.186 | <0.206 |
| -C ₁₂ Aliphatics | | | | | | | <1.93 | <0.196 | 0.206 | <0.153 | 0.321 | <0.102 | <0.109 | <0.172 | <0.136 | <0.176 | <0.2 | <0.127 | <0.222 | <0.205 |
| 9-C ₁₀ Aromatics | | | | | | | <1.93 | <0.196 | 0.747 | <0.153 | 0.786 | <0.102 | <0.109 | <0.172 | <0.136 | <0.176 | <0.2 | <0.127 | <0.222 | < 0.205 |

Notes: For compounds detected at least once above the detection limit, samples reported as not detected (ND) by the laboratory are assumed to have a concentration of one-half of the method detection limit for that sample.

Concentrations entered as < indicate that they were below the detection limit.

Bold print indicates chemicals with no Method 1 standards which are identified as Extractable Petroleum Hydrocarbons (EPH) components. Therefore, EPH (C9-C18 aliphatics) standards are applied.

*MADEP, Background Levels of Polycyclic Aromatic Hydrocarbons and Metals in Soil, *Technical Update*, May 2002.

The maximum concentrations of all metals detected in Site soils with the exception of beryllium and chromium are detected in isolated soils (located >15 feet below the ground surface). NA = Not Available

| ocation: ample Name: | Walpole Pk S. RIZ-1 | Walpole Pk S. RIZ-2 | Walpole Pk S. RIZ-3 | Walpole Pk S. RIZ-4 | Walpole Pk S. RIZ-5 | Walpole Pk S. RIZ-6 | Walpole Pk S. RIZ-7 | Walpole Pk S. RIZ-8 | Walpole Pk S. RIZ-9 | Walpole Pk S. RIZ-10 | | | | | | | | | | | | | |
|--|------------------------|------------------------|------------------------|------------------------|------------------------|------------------------|------------------------|------------------------|------------------------|-------------------------|----------|--------|---------------|---------------|---------------|---------------|---------------|-------------------|-------|----------|----------|----------|---------------------|
| ample Name: ample Depth: | 10'-12' | 15'-17' | 30'-32' | 10'-12' | 1'-3' | 5'-7' | 5'-7' | 812-6 9'-11' | 812-9 9'-11' | 44'-46' | | | | | | | | | | | | | |
| aboratory: | Con-test | Alpha | Alpha | Alpha | Number | Number | Minimum | Maximum | Average | | Soil EPCs | - (malles) | | Method 2 | Method 2 | Method 2 | Concentration |
| aboratory: aboratory I.D.: | 06B06407 | 06B06408 | 06B06409 | 06B06410 | 06B06411 | 06B06412 | 06B06413 | Aipna L018223-01 | Aipria L018223-02 | Aipna L018223-03 | | | Concentration | Concentration | Concentration | Cail | | S (mg/kg) Soil | Soil | Standard | | Standard | In Natural |
| | 15-Feb-06 | | 16-Feb-06 | | 16-Feb-06 | | 16-Feb-06 | | | 6-Dec-07 | | | | | | Soil EPC-1 | Soil EPC-2 | EPC-3 | EPC-4 | S-1/GW-1 | | S-1/GW-3 | In Naturai Soil* |
| ample Date: | | 15-Feb-06 | | 15-Feb-06 | | 16-Feb-06 | | 5-Dec-07 | 6-Dec-07 | | Detected | Sought | Detected | Detected | Detected | | | | | 5-1/GW-1 | 5-1/GW-2 | 5-1/GW-3 | |
| onsultant: | TTR | | | | | | 0-3' | 3'-7' | 8'-15' | >15' | | | | (mg/kg) |
| cetone | <0.059 | < 0.054 | < 0.055 | < 0.047 | 0.059 | 0.11 | <0.075 | 0.0073 | <0.0065 | <0.0055 | 3 | 24 | 0.01 | 0.11 | 0.20 | 0.29 | 0.13 | 0.14 | 0.19 | 6 | 50 | 400 | NA |
| hylbenzene | <0.001 | <0.002 | < 0.002 | <0.001 | 0.001 | <0.002 | <0.002 | < 0.00063 | < 0.00065 | < 0.00055 | 1 | 24 | 0.001 | 0.001 | 0.010 | 0.01 | 0.00 | 0.01 | 0.01 | 40 | 500 | 500 | NA |
| opropylbenzene | <0.001 | < 0.002 | < 0.002 | <0.001 | <0.001 | 0.002 | < 0.002 | < 0.00063 | < 0.00065 | < 0.00055 | 1 | 24 | 0.002 | 0.002 | 0.010 | 0.01 | 0.01 | 0.01 | 0.01 | 1,000 | 1,000 | 1,000 | NA |
| opropyltoluene, p- | < 0.001 | < 0.002 | < 0.002 | < 0.001 | <0.001 | 0.015 | < 0.002 | < 0.00063 | < 0.00065 | < 0.00055 | 1 | 24 | 0.02 | 0.02 | 0.010 | 0.01 | 0.01 | 0.01 | 0.01 | 1,000 | 1,000 | 1,000 | NA |
| ethylene chloride | | | | | | | | | | | | | | | | | | | | | | | |
| lichloromethane) | 0.042 | < 0.011 | 0.030 | < 0.009 | < 0.010 | < 0.011 | < 0.015 | < 0.00063 | < 0.0065 | < 0.0055 | 2 | 24 | 0.03 | 0.04 | 0.10 | 0.14 | 0.05 | 0.07 | 0.10 | 0.1 | 20 | 200 | NA |
| oluene | <0.001 | <0.002 | <0.002 | <0.001 | <0.001 | 0.002 | <0.002 | <0.00094 | <0.00097 | <0.00082 | 1 | 24 | 0.002 | 0.002 | 0.010 | 0.01 | 0.01 | 0.01 | 0.01 | 30 | 500 | 500 | NA |
| rsenic. Total | <6.04 | <5.67 | <5.60 | <5.35 | <5.4 | <6.55 | <5.74 | 1.8 | 4.0 | 3.2 | 5 | 24 | 1.8 | 4.1 | 2.34 | 2.0 | 2.6 | 2.7 | 2.2 | 20 | 20 | 20 | 20 |
| arium. Total | 29.0 | 42.1 | 16.9 | 18.5 | 16.4 | 23.9 | 21.8 | 18.0 | 32.0 | 37.0 | 24 | 24 | 9.7 | 42.7 | 23.4 | 16.1 | 26.9 | 26.3 | 28.1 | 1,000 | 1,000 | 1,000 | 50 |
| eryllium, Total | 0.45 | 0.87 | 0.42 | <0.27 | <0.27 | 0.37 | 0.36 | 0.40 | 0.83 | 0.88 | 8 | 10 | 0.36 | 0.88 | 0.49 | 0.14 | 0.37 | 0.45 | 0.72 | 100 | 100 | 100 | 0.4 |
| hromium. Total | 14.4 | 6.4 | 8.2 | 6.7 | 6.5 | 10.7 | 5.6 | 8.3 | 6.2 | 6.1 | 24 | 24 | 3.9 | 30.7 | 9.2 | 9.1 | 7.28 | 7.5 | 12.5 | 30 | 30 | 30 | 30 |
| ead, Total | 6.8 | 6.3 | 4.8 | 6.8 | 10.6 | 14.5 | 12.1 | 3.8 | 19.0 | 18.0 | 23 | 30 | 3.1 | 90.1 | 10.4 | 4.7 | 10.3 | 16.2 | 9.7 | 300 | 300 | 300 | 100 |
| ickel, Total | 7.2 | 3.8 | 6.7 | 5.64 | 4.7 | 6.4 | 4.7 | 6.3 | 5.0 | 3.7 | 10 | 10 | 3.7 | 7.2 | 5.4 | 4.7 | 5.5 | 6.0 | 4.7 | 20 | 20 | 20 | 20 |
| anadium | 13.0 | 7.7 | 12.2 | 9.8 | 12.5 | 22.2 | 10.6 | 16.0 | 12.0 | 7.7 | 10 | 10 | 7.7 | 22.2 | 12.4 | 12.5 | 16.4 | 12.7 | 9.2 | 600 | 600 | 600 | 30 |
| nc, Total | 32.3 | 41.4 | 27.4 | 23.2 | 20.3 | 24.5 | 24.9 | 34.0 | 45.0 | 70.0 | 10 | 10 | 20.3 | 70.0 | 34.3 | 20.3 | 24.7 | 33.6 | 46.3 | 2,500 | 2,500 | 2,500 | 100 |
| ercury, Total | < 0.011 | <0.006 | < 0.005 | <0.011 | 0.016 | 0.031 | 0.029 | <0.08 | <0.08 | <0.09 | 3 | 24 | 0.02 | 0.03 | 0.06 | 0.08 | 0.05 | 0.05 | 0.06 | 20 | 20 | 20 | 0.3 |
| ₉ -C ₁₂ Aliphatics | | | | | | | | | | | 2 | 14 | 0.21 | 0.32 | 0.17 | 0.26 | 0.06 | 0.08 | 0.09 | 1,000 | 1,000 | 1,000 | NA |
| ₉ -C ₁₀ Aromatics | | | | | | | | | | | 2 | 14 | 0.75 | 0.79 | 0.24 | 0.40 | 0.06 | 0.08 | 0.09 | 100 | 100 | 100 | NA |

Notes: For compounds detected at least once above the detection limit, samples reported as not detected (ND) by the laboratory

are assumed to have a concentration of one-half of the method detection limit for that sample.

Concentrations entered as < indicate that they were below the detection limit.

Bold print indicates chemicals with no Method 1 standards which are identified as Extractable Petroleum Hydrocarbons (EPH) components. Therefore, EPH (C9-C18 aliphatics) standards are applied.

*MADEP, Background Levels of Polycyclic Aromatic Hydrocarbons and Metals in Soil, *Technical Update*, May 2002.

The maximum concentrations of all metals detected in Site soils with the exception of beryllium and chromium are less than "natural" soil background. The maximum concentrations of beryllium and chromium are detected in isolated soils (located >15 feet below the ground surface). NA = Not Available

| Table 3 | Positive Groundwater Ar | alytical Data (µg/L) - Walpole Park S | outh, Walpole, Massachusetts |
|-----------|-------------------------|---------------------------------------|------------------------------|
| Location: | Walpole Pk S. Walpole | Pk S. Walpole Pk S. Walpole Pk S | S. Walpole Pk S |

| Location: | | | | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | | Walpole Pk S. | | Walpole Pk S. | |
|-------------------------|----------|----------|----------|---------------|---------------|---------------|---------------|--------------|---------------|---------------|---------------|---------------|---------------|---------------|--------------|---------------|---------------|---------------|---------------|---------------|---------------|--------------|
| Sample Name: | | | | MW-2-042808 | MW-2-GW | MW-2 | MW-2 | MW-2 | MW-3 | MW-3-042808 | MW-3-GW | MW-3 | MW-3 | MW-3 | MW-3 | MW-9 | MW-9-051408 | MW-9-GW | MW-9 | MW-9 | MW-9 | MW-9 |
| Laboratory: | Method 2 | Method 2 | Method 2 | Alpha | Spectrum | Alpha | Alpha | Well Average | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average |
| Laboratory I.D.: | Standard | Standard | Standard | L0806023-07 | SA87371-05 | L0907670-07 | L1008812-03 | | L0718979-07 | L0806023-03 | SA87371-07 | L0908197-01 | L0918777-07 | L1008812-01 | | L0718979-05 | L0806993-02 | SA87371-04 | L0907670-04 | L0918777-01 | L1008812-08 | |
| Sample Date: | GW-1 | GW-2 | GW-3 | 28-Apr-08 | 11-Nov-08 | 10-Jun-09 | 11-Jun-10 | | 20-Dec-07 | 28-Apr-08 | 11-Nov-08 | 8-Jun-09 | 28-Dec-09 | 11-Jun-10 | | 20-Dec-07 | 14-May-08 | 11-Nov-08 | 10-Jun-09 | 21-Dec-09 | 11-Jun-10 | 1 |
| Consultant: | | | | TTR | TTR | TTR | TTR | | TTR | TTR | TTR | TTR | TTR | TTR | | TTR | TTR | TTR | TTR | TTR | TTR | |
| Chloroform | 70 | 50 | 20,000 | 2.2 | <1.0 | 1.0 | 1.2 | 1.2 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.75 | 0.71 | 2.2 | 0.75 | <0.5 | <0.5 | 0.8 |
| Methyl tert-butyl ether | 70 | 50,000 | 50,000 | <0.50 | <1.0 | <0.50 | <0.5 | 0.3 | 1.7 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.5 | 5.1 | 7.9 | 2.2 | <0.50 | 0.77 | <0.5 | 2.8 |
| Toluene | 1,000 | 6,000 | 50,000 | <0.50 | <1.0 | <0.50 | <0.5 | 0.3 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 |
| Barium, Dissolved | 2,000 | NA | 50,000 | 31.0 | 120 | 70 | 70 | 75.0 | 15.2 | 10.0 | 11.4 | 24 | -10 | 36 | 16.4 | 7.0 | 13.0 | 18.6 | 20 | 56 | 64 | 31.3 |
| Lead, Dissolved | 15 | NΑ | 10 | <10.0 | <7.5 | <10 | /U /10 | 15.0 4.7 | -2.0 | <10.0 | 8.8 | Z1 Z10 | <10 | -10 | 5.0 | <2.0 | <10.0 | <7.5 | 29 -10 | <10 | <10 | A 1 |
| Nickel, Dissolved | 100 | NA NA | 200 | <25.0 | <5.0 | <25 | <25 | 10.0 | <2.0 | <25.0 | <5.0 | <25 | <25 | <25 | 8.9 | <2.0 | <10.0 | <5.0 | <25 | <25 | <25 | 8.2 |
| Thallium, Dissolved | 2 | NA. | 400 | <2.0 | <5.0 | <2.0 | <2 | ND | <2.0 | <2.0 | <5.0 | <2.0 | <2 | <2 | ND | <2.0 | | <5.0 | <2.0 | <2 | <2 | ND |
| Zinc, Dissolved | 5,000 | NA | 900 | <50.0 | 28.0 | <50 | <50 | 25.8 | <20.0 | <50.0 | 34.5 | <50 | <50 | <50 | 24.1 | 25.9 | | 34.7 | <50 | 52 | 51 | 37.7 |
| | | | | | | | | | | | | | | | | | | | | | ŀ | 1 |

For compounds detected at least once above the detection limit, samples reported as not detected (ND) by the laboratory are assumed to have a concentration of one-half of the method detection limit for that sample. Concentrations entered as < indicate that they were below the detection limit.

NA = Not Applicable; ND = Not Detected

P:\Pre-FY2008\12700000\12700058\Risk Assessment\RA_RAO_Jan 2011\12700058-Walpole Soil & GW data_2010.xlsGW_DATA-2010 RA

| ocation: | | | | Walpole Pk S. | | Walpole Pk S. | | Walpole Pk S. | Walpole Pk S. | | Walpole Pk S. | Walpole Pk S. | |
|---------------------|----------|----------|----------|---------------|---------------|---------------|---------------|---------------|---------------|--------------|---------------|---------------|---------------|---------------|---------------|---------------|--------------|---------------|---------------|--------------|---------------|---------------|---------------|---------------|------------------|---------------|--------------|
| mple Name: | | | | RIZ-3 | RIZ-3-051408 | RIZ-3-GW | RIZ-3 | RIZ-3 | RIZ-3 | RIZ-3 | RIZ-8 | RIZ-8-042808 | RIZ-8-GW | RIZ-8 | RIZ-8 | RIZ-8 | RIZ-8 | RIZ-8S-042808 | RIZ-8S | RIZ-8S | RIZ-9 | RIZ-9-042808 | RIZ-9-GW | RIZ-9 | RIZ-9 | RIZ-9 | RIZ-9 |
| oratory: | Method 2 | Method 2 | Method 2 | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average | Alpha | ALPHA | Well Average | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average |
| oratory I.D.: | Standard | | Standard | L0718979-06 | L0806993-01 | SA87371-03 | L0907670-06 | L0918777-04 | L1008812-02 | | L0718979-01 | L0806023-04 | SA87371-02 | L0907670-02 | L0918777-02 | L1008812-07 | | L0806023-05 | L0907670-03 | | L0718979-03 | L0806023-02 | SA87371-06 | L0907670-08 | L0918777-05 | L1008812-05 | |
| mple Date: | GW-1 | GW-2 | GW-3 | 20-Dec-07 | 14-May-08 | 11-Nov-08 | 10-Jun-09 | 28-Dec-09 | 11-Jun-10 | | 19-Dec-07 | 28-Apr-08 | 11-Nov-08 | 10-Jun-09 | 21-Dec-09 | 11-Jun-10 | | 28-Apr-08 | 10-Jun-09 | | 19-Dec-07 | 28-Apr-08 | 11-Nov-08 | 10-Jun-09 | 28-Dec-09 | 11-Jun-10 | |
| nsultant: | | | | TTR | TTR | TTR | TTR | TTR | TTR | | TTR | TTR | TTR | TTR | TTR | TTR | | TTR | TTR | | TTR | TTR | TTR | TTR | TTR | TTR | |
| | 70 | 50 | 00.000 | 0.75 | 0.50 | 4.0 | 0.50 | 0.5 | ٥٠ | 0.0 | 0.75 | 0.50 | 1.0 | 0.50 | ٥٠ | 0.5 | 0.0 | 0.50 | 0.50 | 0.0 | 0.75 | 0.50 | | 0.50 | 2.5 | ٥٠ | 2.0 |
| roform | 70 | 50 | 20,000 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.50 | <0.50 | 0.3 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 |
| yl tert-butyl ether | 70 | 50,000 | 50,000 | <1.0 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <1.0 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.50 | <0.50 | 0.3 | <1.0 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 |
| ene | 1,000 | 6,000 | 50,000 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <5 | 0.7 | <0.50 | 0.71 | 0.5 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 |
| um, Dissolved | 2,000 | NA | 50,000 | 25.6 | 44.0 | 86.2 | 13 | 177 | 161 | 84.5 | 50.8 | 25.0 | 27.6 | 26 | 31 | <10 | 27.6 | 54.0 | 51 | 52.5 | 15.3 | 21.0 | 14.8 | 15 | <10 | <10 | 12.7 |
| l, Dissolved | 15 | NA | 10 | <2.0 | <10.0 | <7.5 | <10 | <10 | <10 | 4.1 | <2.0 | <10.0 | <7.5 | <10 | <10 | <10 | 4.1 | <10.0 | <10 | 5.0 | <2.0 | <10.0 | <7.5 | <10 | <10 | <10 | 4.1 |
| el, Dissolved | 100 | NA | 200 | <2.0 | | <5.0 | <25 | <25 | <25 | 8.2 | 4.8 | <25.0 | <5.0 | <25 | <25 | <25 | 9.6 | <25.0 | <25 | 12.5 | <2.0 | <25.0 | <5.0 | <25 | <25 | <25 | 8.9 |
| lium, Dissolved | 2 | NA | 400 | <2.0 | | <5.0 | <2.0 | <2 | <2 | ND | <2.0 | <2.0 | <5.0 | <2.0 | <2 | <2 | ND | <2.0 | <2.0 | ND | <2.0 | <2.0 | <5.0 | <2.0 | <2 | <2 | ND |
| c, Dissolved | 5,000 | NA | 900 | <20.0 | | 21.0 | ~50 | <50 | <50 | 21.2 | <20.0 | <50.0 | 26.4 | <50 | <50 | <50 | 22.7 | <50.0 | <50 | 25.0 | <20.0 | <50.0 | 20.0 | <50 | ~ 5 0 | <50 | 21.7 |

For compounds detected at least once above the detection limit, samples reported as not detected (ND) by the laboratory are assumed to have a concentration of one-half of the method detection limit for that sample.

Concentrations entered as < indicate that they were below the detection limit.

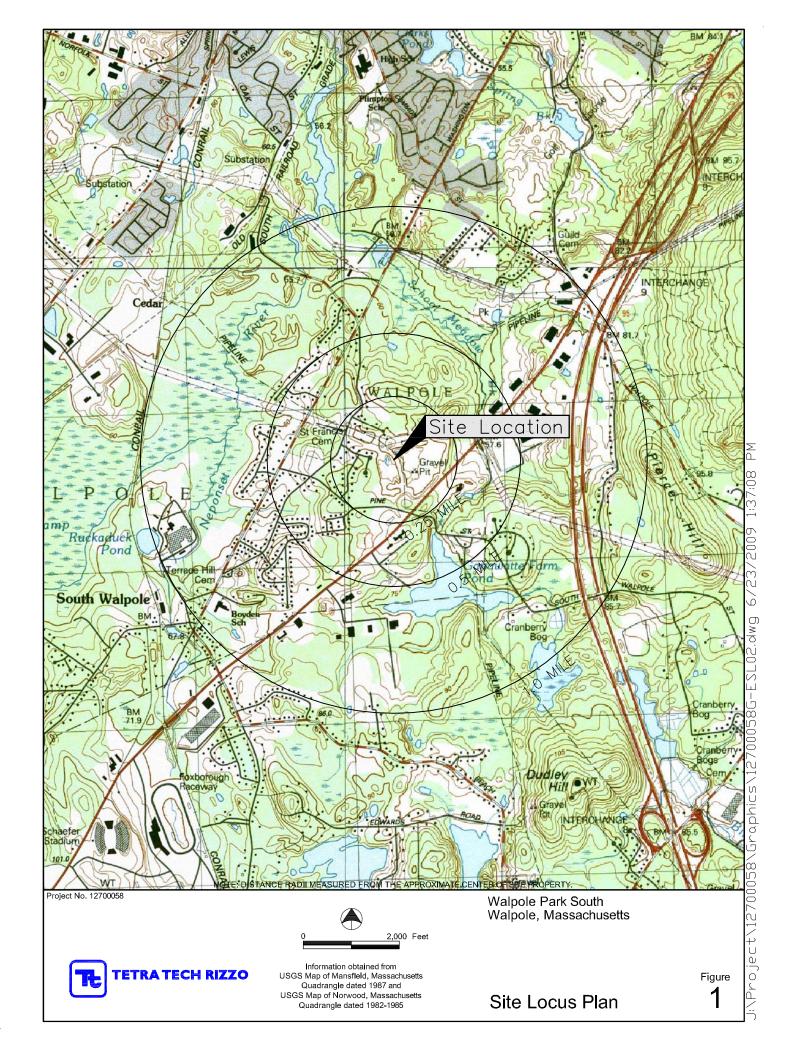
NA = Not Applicable; ND = Not Detected

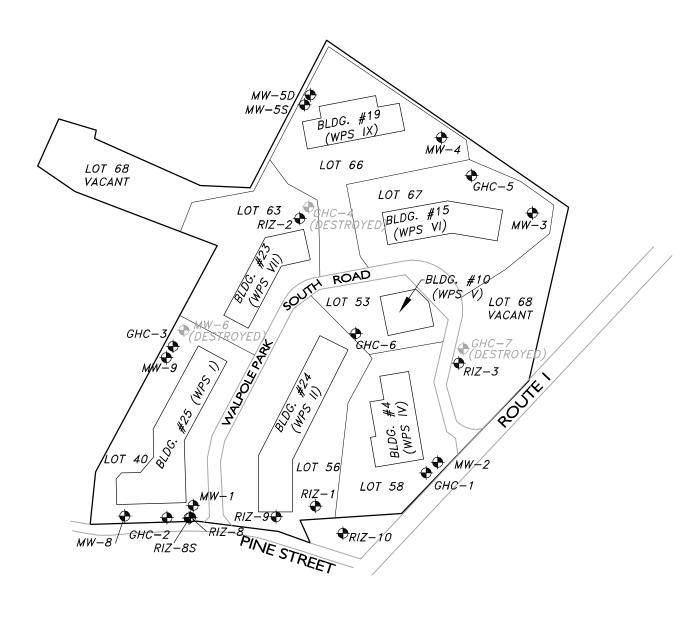
| Table 3 | | | | Positive Grounds | water Analytical Da | ata (µg/L) - Walpo | le Park South, Wa | alpole, Massachu | setts | | | | | | | | | | Groundwate | r Summary S | Statistics (µg/L) | |
|-------------------------|----------|----------|----------|------------------|---------------------|--------------------|-------------------|------------------|---------------|---------------|--------------|---------------|---------------|---------------|---------------|---------------|---------------|--------------|------------|-------------|-------------------|---------------|
| Location: | | | | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | Walpole Pk S. | | Walpole Pk S. | | | | | |
| Sample Name: | | | l I | RIZ-10 | RIZ-10-042808 | RIZ-10-GW | RIZ-10 | RIZ-10 | RIZ-10 | RIZ-10 | RIZ-10 | GHC-6 | GHC-6-042808 | GHC-6-GW | GHC-6 | GHC-6 | GHC-6 | GHC-6 | | | | |
| Laboratory: | Method 2 | Method 2 | Method 2 | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Alpha | Well Average | Alpha | Alpha | Spectrum | Alpha | Alpha | Alpha | Well Average | Number | Number | Minimum | Maximum |
| Laboratory I.D.: | Standard | Standard | Standard | L0718979-02 | L0806023-01 | SA87371-01 | L0818397-01 | L0907670-01 | L0918777-06 | L1008812-06 | _ | L0718979-04 | L0806023-06 | SA87371-08 | L0907670-05 | L0918777-03 | L1008812-04 | _ | of Times | of Times | Concentration | Concentration |
| Sample Date: | GW-1 | GW-2 | GW-3 | 19-Dec-07 | 28-Apr-08 | 11-Nov-08 | 11-Dec-08 | 10-Jun-09 | 28-Dec-09 | 11-Jun-10 | | 19-Dec-07 | 28-Apr-08 | 11-Nov-08 | 10-Jun-09 | 28-Dec-09 | 11-Jun-10 | | Detected | Sought | Detected | Detected |
| Consultant: | | | l I | TTR | TTR | TTR | TTR | TTR | TTR | TTR | | TTR | TTR | TTR | TTR | TTR | TTR | | | - | | |
| | | | | | | | | | | | | | | | | | | | | | | |
| Chloroform | 70 | 50 | 20,000 | < 0.75 | < 0.50 | <1.0 | | < 0.50 | < 0.5 | <0.5 | 0.3 | < 0.75 | < 0.50 | <1.0 | < 0.50 | <0.5 | <0.5 | 0.3 | 6 | 48 | 0.7 | 2.2 |
| Methyl tert-butyl ether | 70 | 50,000 | 50,000 | 1.2 | < 0.50 | <1.0 | | < 0.50 | < 0.5 | < 0.5 | 0.5 | <1.0 | < 0.50 | <1.0 | < 0.50 | < 0.5 | < 0.5 | 0.3 | 6 | 48 | 0.8 | 7.9 |
| Toluene | 1,000 | 6,000 | 50,000 | <0.75 | 0.73 | <1.0 | | <0.50 | <0.5 | <0.5 | 0.4 | <0.75 | <0.50 | <1.0 | <0.50 | <0.5 | <0.5 | 0.3 | 2 | 48 | 0.7 | 0.7 |
| Barium, Dissolved | 2,000 | NA | 50,000 | 95.8 | 62.0 | 88.4 | | 148 | 99 | 107.0 | 100.0 | 45.9 | 59.0 | 36.8 | 66 | 39 | 63 | 51.6 | 44 | 48 | 7.0 | 177.0 |
| Lead, Dissolved | 15 | NA | 10 | <2.0 | <10.0 | <7.5 | | <10 | <10 | <10 | 4.1 | <2.0 | <10.0 | <7.5 | <10 | <10 | <10 | 4.1 | 1 | 48 | 8.8 | 8.8 |
| Nickel, Dissolved | 100 | NΔ | 200 | 7.9 | <25.0 | <5.0 | | <25 | <25 | <25 | 10.1 | <2.0 | <25.0 | <5.0 | <25 | -25 | <25 | 8.0 | 2 | 46 | 4.8 | 7.9 |
| Thallium, Dissolved | 2 | NΔ | 400 | <2.0 | <2.0 | 11.6 | <2.0 | <2 | -2 | -20 | ND* | <2.0 | <2.0 | <5.0 | <2.0 | -2 | <2 | ND | 1 | 47 | 11.6 | 11.6 |
| Zinc, Dissolved | 5,000 | NA | 900 | 21.6 | <50.0 | 36.3 | ~2.0 | <50 | <50 | <50 | 26.3 | <20.0 | <50.0 | 21.6 | <50 | -50 | <50 | 21.9 | 12 | 46 | 20.0 | 52.0 |
| Ziric, Dissolved | 3,000 | INA | 300 | 21.0 | < 30.0 | 30.3 | | 230 | 230 | 230 | 20.3 | ₹20.0 | <50.0 | 21.0 | 250 | 230 | <50 | 21.9 | 12 | 40 | 20.0 | 32.0 |

For compounds detected at least once above the detection limit, samples reported as not detected (ND) by the laboratory are assumed to have a concentration of one-half of the method detection limit for that sample.

Concentrations entered as < indicate that they were below the detection limit.

NA = Not Applicable; ND = Not Detected; * Thallium detected in RIZ-10 during the November 2008 round of sampling is considered anomalous and therefore, it is not considered to be a COC at the Site. See text for details.

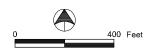




LEGEND

- LOT BOUNDARIES

12700058P-ESP02

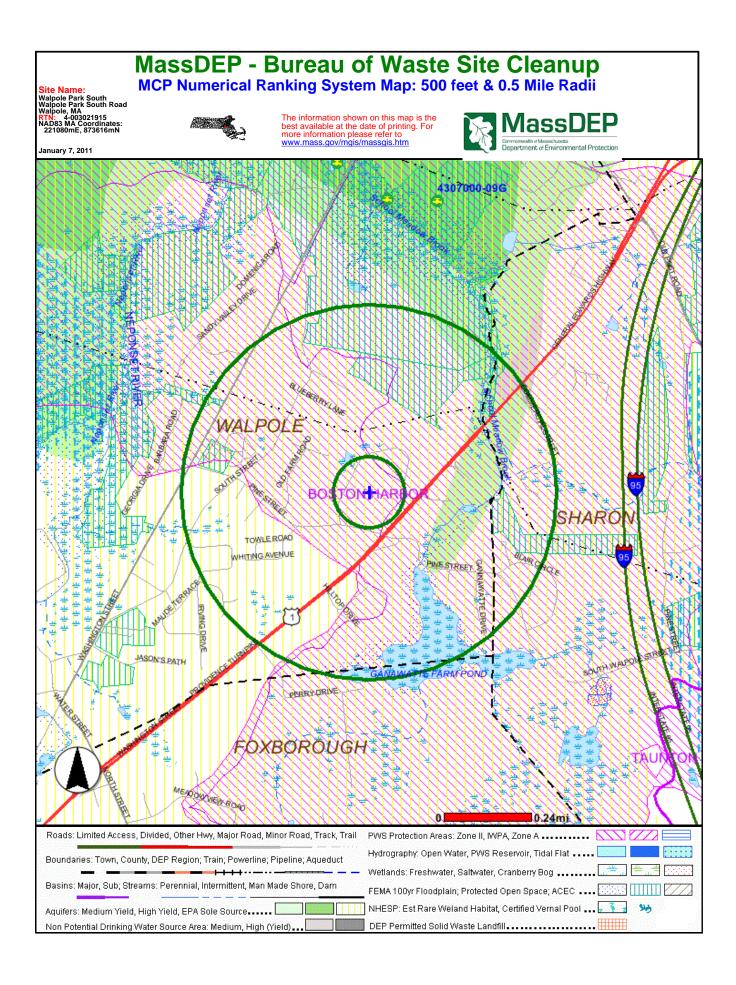


Walpole Park South Walpole, Massachusetts

Site Plan with Monitoring Well Locations



Site Plan by GeoHydroCycle, Inc. Dated 5/14/04 Map Output Page 1 of 1



Appendix A LSP Statement of Limitations and Conditions

Statement of Limitations and Conditions

Attachment to Opinion of Massachusetts Licensed Site Professional

Tetra Tech Rizzo

Name of Licensed Site Professional: Raymond C. Johnson

LSP Registration Number: 6118

Date of Opinion: January 10, 2011

Client to Whom Opinion was Walpole Park South Trust

Rendered:

Response Tracking No./Site No.: 4-3021915

This Statement of Limitations and Conditions is an integral part of, and is incorporated by reference into, the Opinion of Massachusetts Licensed Site Professional referenced above.

Limitations

1. Purpose of Opinion

- A. This Opinion is being provided in compliance with the requirements set forth in the Massachusetts Contingency Plan ("MCP"), 310 CMR 40.0000 et seq. Specifically, the LSP has prepared this Opinion at the request of the Client identified above as part of a Class A-2 Response Action Outcome Statement/Phase V Completion Statement. This stated purpose has been a significant factor in determining the scope and level of services required to render this Opinion.
- B. Should the purpose for which this Opinion is to be used change, this Opinion shall no longer be valid.

2. General

A. This Opinion was prepared for the sole and exclusive use of the Client, subject to the provisions of the MCP. No other party is entitled to rely in any way on the conclusions, observations, specifications, or data contained herein without the express written consent of Tetra Tech Rizzo and the LSP who rendered this opinion. Any use of this Opinion by anyone other than Client, or any use of this Opinion by Client or others for any purpose other than the stated purpose set forth above, without the LSP's review and the written authorization of Tetra Tech Rizzo and the LSP, shall be at the user's sole risk, and neither Tetra Tech Rizzo nor the LSP shall have any liability or responsibility therefor.

B. This Opinion was prepared pursuant to an Agreement between Tetra Tech Rizzo and the Client referenced above which defines the scope of work and sets out agreements regarding waivers of consequential damages, limitations on liability, and other important conditions and restrictions pursuant to which the Opinion is rendered. All uses of the Opinion are subject to and deemed acceptance of the conditions and restrictions contained in such Agreement. A copy of the Agreement or relevant excerpts from the Agreement will be made available upon requests to any authorized person seeking to use the Opinion.

3. Scope of Services

The observations and conclusions described in this Opinion are based solely on the Services provided pursuant to the Agreement with the Client and any approved additional services authorized by Client. Without limitation of any other applicable limitations or conditions, neither Tetra Tech Rizzo nor the LSP shall be liable for the existence of any condition, the discovery of which would have required the performance of services not authorized under the Agreement. To the best of the knowledge and belief of Tetra Tech Rizzo and the LSP who signed this Opinion, no inquiry of an attorney-at-law having being made, no laws, regulations, orders, permits or approvals are applicable to the response actions to which this opinion relates except, if and to the extent applicable, M.G.L. c. 21A, Sections 19-19J, 309 CMR, M.G.L. c. 21 E and 310 CMR 40.0000. Accordingly, this opinion is not intended to and does not address compliance with any other laws, regulation, orders, permits or approvals.

4. Changed Circumstances

The passage of time may result in changes in technology, economic conditions or regulatory standards, manifestations of latent conditions, or the occurrence of future events which would render this Opinion inaccurate or otherwise inapplicable. Neither Tetra Tech Rizzo nor the LSP shall be liable or responsible for the consequences of any such changed circumstances or conditions on the accuracy of this Opinion. In addition, under no circumstances shall the Client nor any other person or entity rely on the information or conclusions contained in this Opinion after six months from its date of submission without the express written consent of Tetra Tech Rizzo and the LSP. Reliance on the Opinion after such period of time shall be at the user's sole risk.

- **5.** Should Tetra Tech Rizzo or the LSP be required or requested to review or authorize others to use this Opinion after its date of submission, Tetra Tech Rizzo shall be entitled to additional compensation at then existing rates or such other terms as may be agreed upon between Tetra Tech Rizzo and the Client. Nothing herein contained shall be deemed to require Tetra Tech Rizzo or the LSP to undertake any such review or authorize others to use this Opinion.
- **6.** The conclusions stated in this Opinion are based upon:

Statement of Limitations and Conditions Attachment to Opinion of Massachusetts Licensed Site Professional

- Visual inspection of existing physical conditions;
- Review and interpretation of site history and site usage information which was made available or obtained within the scope of work authorized by the Client;
- Information provided by the Client;
- Information and/or analyses for designated substances or parameters provided by an independent testing service or laboratory on a limited number of samples; and
- A limited number of subsurface explorations made on dates indicated in documentation supporting this Opinion;

The information upon which the LSP has relied and presumed accurate, and upon which the LSP is entitled to reasonably rely. The LSP was not authorized and did not attempt to independently verify the accuracy or completeness of information or materials received from the Client and/or from laboratories and other third parties during the performance of its services. Neither Tetra Tech Rizzo nor the LSP shall be liable for any condition, information, or conclusion, the discovery of which required information not available to the LSP or for independent investigation of information provided to the LSP by the Client and/or independent third parties.

7. This Opinion is rendered for the limited purpose stated above, and is not and should not be deemed to be an opinion concerning the compliance of any past or present owner or operator of the site with any federal, state or local law or regulation. No warranty or guarantee, whether express or implied, is made by this opinion, and any implied warranties of merchantability or fitness for a particular purpose are expressly disclaimed. Without limiting the generality of the foregoing, no warranty or guarantee is made that all contamination at a site or sources or contamination has been detected or identified, that any action or recommended action will achieve all of its objectives, or that this Opinion or any action as to which this Opinion relates will be upheld by any audit conducted by the DEP or any other party.

Appendix A: Limitations

- 1. The observations described in this report were made under the conditions stated therein. The conclusions presented in the report were based solely upon the services described therein, and not on scientific tasks or procedures beyond the scope of described services or the time and budgetary constraints imposed by the CLIENT. The work described in this report was carried out in accordance with the Terms and Conditions in our contract.
- In preparing this report, ENGINEER has relied on certain information provided by state and local officials and other parties referenced therein, and on information contained in the files of state and/or local agencies available to ENGINEER at the time of the site assessment. Although there may have been some degree of overlap in the information provided by these various sources, ENGINEER did not attempt to independently verify the accuracy or completeness of all information reviewed or received during the course of this site assessment.
- 3. Observations were made of the Site and of structures on the Site as indicated within the report. Where access to portions of the Site or to structures on the Site was unavailable or limited, ENGINEER renders no opinion as to the presence of hazardous materials or oil, or to the presence of indirect evidence relating to hazardous material or oil, in that portion of the Site or structure. In addition, ENGINEER renders no opinion as to the presence of hazardous material or oil, or the presence of indirect evidence relating to hazardous material or oil, where direct observation of the interior walls, floor, or ceiling of a structure on a Site was obstructed by objects or coverings on or over these surfaces.
- 4. ENGINEER did not perform testing or analyses to determine the presence or concentration of asbestos at the Site or in the environment at the Site.
- 5. It is ENGINEER's understanding that the purpose of this report is to assess the physical characteristics of the subject Site with respect to the presence on the Site of hazardous material or oil. This stated purpose has been a significant factor in determining the scope and level of services provided for in the Agreement. Should the purpose for which the Report is to be used or the proposed use of the site(s) change, this Report is no longer valid and use of this Report by CLIENT or others without ENGINEER's review and written authorization shall be at the user's sole risk. Should ENGINEER be required to review the Report after its date of submission, ENGINEER shall be entitled to additional compensation at then existing rates or such other terms as agreed between ENGINEER and the CLIENT.
- 6. The conclusions and recommendations contained in this report are based in part, where noted, upon the data obtained from a limited number of soil samples obtained from widely spaced subsurface explorations. The nature and extent of variations between these explorations may not become evident until further exploration. If variations or other latent conditions then appear evident, it will be necessary to reevaluate the conclusions and recommendations of this report.
- 7. Any water level readings made in test pits, borings, and/or observation wells were made at the times and under the conditions stated on the report. However, it must be noted that fluctuations in the level of groundwater may occur due to variations in rainfall and other factors different from those prevailing at the time measurements were made.

- 8. Except as noted within the text of the report, no quantitative laboratory testing was performed as part of the site assessment. Where such analyses have been conducted by an outside laboratory, ENGINEER has relied upon the data provided and has not conducted an independent evaluation of the reliability of these data.
- 9. The conclusions and recommendations contained in this report are based in part, where noted, upon various types of chemical data and are contingent upon their validity. These data have been reviewed and interpretations made in the report. As indicated within the report, some of these data may be preliminary screening level data and should be confirmed with quantitative analyses if more specific information is necessary. Moreover, it should be noted that variations in the types and concentrations of contaminants and variations in their flow paths may occur due to seasonal water table fluctuations, past disposal practices, the passage of time, and other factors. Should additional chemical data become available in the future, these data should be reviewed, and the conclusions and recommendations presented herein modified accordingly.
- 10. Chemical analyses have been performed for specific constituents during the course of this site assessment, as described in the text. However, it should be noted that additional chemical constituents not searched for during the current study may be present in soil and/or groundwater at the Site.
- 11. This Report was prepared for the exclusive use of the CLIENT. No other party is entitled to rely on the conclusions, observations, specifications, or data contained therein without the express written consent of ENGINEER.
- 12. The observations and conclusions described in this Report are based solely on the Scope of Services provided pursuant to the Agreement. ENGINEER has not performed any additional observations, investigations, studies, or testing not specifically stated therein. ENGINEER shall not be liable for the existence of any condition, the discovery of which required the performance of services not authorized under the Agreement.
- 13. The passage of time may result in significant changes in technology, economic conditions, or site variations that would render the Report inaccurate. Accordingly, neither the CLIENT, nor any other party, shall rely on the information or conclusions contained in this Report after six months from its date of submission without the express written consent of ENGINEER. Reliance on the Report after such period of time shall be at the user's sole risk. Should ENGINEER be required to review the Report after six months from its date of submission, ENGINEER shall be entitled to additional compensation at then existing rates or such other terms as may be agreed upon between ENGINEER and the CLIENT.
- 14. ENGINEER has endeavored to perform its services based upon engineering practices accepted at the time they were performed. ENGINEER makes no other representations, express or implied, regarding the information, data, analysis, calculations, and conclusions contained herein.
- 15. The services provided by ENGINEER do not include legal advice. Legal counsel should be consulted regarding interpretation of applicable and relevant federal, state, and local statutes and regulations and other legal matters.

Appendix B

Data Usability Assessment Documentation

| Appendix B - Data Usability As | sessment Summary Table | 1 | , | | , | | , , | | | | | , | , | / . | , | | |
|--------------------------------|-------------------------------------|-----------------|---------|-------|-----------------------|-----|--------------------|-------------------|-----------|-------------------|-----------|---------------------------------|-----------------|-------------|-------|------------------|---|
| Sample ID or Series | Parameters | Date Sampled | los los | Sedim | Site Characterization | FPS | Hazard Elimination | CAM Compliant (X) | Sample Ro | Method QA/QC Proc | QA/QC Ro. | Achieved? (Y/N) Trip Blanks OV. | Field Duplicate | Relative D. | (RPD) | Limite Reporting | Data Qualifications |
| RIZ-8 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | x | х | | Υ | Υ | Y | Y | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc oritera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are "difficult analytes". None of these compounds are COCs at the Site. Metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-10 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | х | х | | Υ | Υ | Y | Υ | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc critera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are "difficult analytes". None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-9 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | х | х | | Y | Y | Y | Y | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc critera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are "difficult analytes". None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| GHC-6 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | x | х | | Υ | Y | Y | Y | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc oritera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are 'difficult analytes'. None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| MW-9 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | x | х | | Υ | Υ | Υ | Y | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc oritera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are 'difficult analytes'. None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-3 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | X | | х | х | | Υ | Y | Y | Υ | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc critera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are 'difficult analytes'. None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| MW-3 | VOCs by 8260, EPH, MCP 14 Metals | 19-Dec-07 | х | | x | х | | Υ | Υ | Υ | Y | NA | NA | | Y | | LCS/LCSD % recoveries for Dichlorodifluoromethand, 1,4-Dioxane and Acetone are outside of the individual acceptanc oritera for the compounds, but within overall method allowances. The laboratory report indicates that these three compounds are 'difficult analytes'. None of these compounds are COCs at the Site. Metals samples were diluted due the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-10 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | Х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Υ | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-9 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | Х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| MW-3 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Υ | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-8 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-8S | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| GHC-6 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| MW-2 | VOCs by 524.2, MCP 14 Metals | 28-Apr-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | Some metals samples were diluted due to the presence of non-target analytes. RDLs are at or below applicable MCP Method 1 standards. |
| RIZ-10 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | х | | Х | х | | Υ | Υ | Υ | Υ | Υ | NA | | Υ | | mound i dundado. |
| RIZ-8 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| RIZ-3 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| MW-9 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| MW-2 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| RIZ-9 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| MW-3 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| GHC-6 | VOCs by 624, MCP 14 Metals | 11-Nov-08 | Х | | х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Y | | |
| RIZ-10 | Thallium | 11-Dec-08 | Х | | х | х | | Υ | Υ | Y | Υ | NA | NA | | Y | | Sample received by laboratory beyoned recommended holding time for laboratory filtration. Not considered to be significant based on absence of turbidity or observed sediment in sample. Sample was diluted because of non-target analytes, but applicable MCP reporting limits were achieved. |
| RIZ-10 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | Х | | Х | х | | Υ | Υ | Υ | Υ | NA | NA | | Y | | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |

| Sample ID or Series | Parameters | Date Sampled | / | Soil | Secient | Site Characte | Tation | Hazard Elimi | CAM Complies | Sample R. | Method QA/OC | Followed? (YM) QA/QC Ro. | Achieved? (Y/N) Trip Blanks OV. | Field Duplicate C. | Relative Percent | Appropriate S | Data Qualifications |
|---------------------|--|-----------------|---|----------|---------|------------------|--------------|--------------|--------------|-----------|--------------|--------------------------|--|--------------------|------------------|---------------|--|
| RIZ-8 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | х | х | | Υ | Υ | Y | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| RIZ-8S | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | х | Х | | Υ | Υ | Y | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| MW-9 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | х | Х | | Υ | Υ | Υ | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| GHC-6 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | х | Х | | Υ | Υ | Y | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| RIZ-3 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | х | Х | | Υ | Υ | Y | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| MW-2 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | Х | | Х | Х | | Υ | Υ | Υ | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| RIZ-9 | VOCs by 524.2, MCP 14 Metals | 10-Jun-09 | | х | | Х | Х | | Υ | Υ | Y | Υ | NA | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| MW-9 | VOCs by 524.2, MCP 14 Metals | 21-Dec-09 | | х | | Х | Х | | Υ | Υ | Υ | Υ | Υ | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| RIZ-8 | VOCs by 524.2, MCP 14 Metals | 21-Dec-09 | | Х | | Х | Х | | Υ | Υ | Y | Υ | Υ | NA | | Υ | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe |
| GHC-6 | VOCs by 524.2, MCP 14 Metals | 28-Dec-09 | | Х | | Х | Х | | Υ | Υ | Y | Y | Υ | NA | | Υ | analytes. The applicable MCP reporting limits were achieved. Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe analytes. The applicable MCP reporting limits were achieved. |
| RIZ-3 | VOCs by 524.2, MCP 14 Metals | 28-Dec-09 | | Х | | Х | Х | | Υ | Y | Y | Y | Υ | NA | | Y | Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe |
| RIZ-9 | VOCs by 524.2, MCP 14 | 28-Dec-09 | | Х | | Х | Х | | Y | Y | Y | Y | Υ | NA | | Y | analytes. The applicable MCP reporting limits were achieved. Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe |
| RIZ-10 | Metals VOCs by 524.2, MCP 14 | 28-Dec-09 | | Х | | х | Х | | Y | Y | Y | Y | Υ | NA | | Y | analytes. The applicable MCP reporting limits were achieved. Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe |
| MW-3 | Metals VOCs by 524.2, MCP 14 | 28-Dec-09 | | X | | Х | Х | | Y | Y | Y | Y | Y | NA | | Y | analytes. The applicable MCP reporting limits were achieved. Metals samples have high detection limits for Antimony and Thallium due to dilutions required by presence of non-targe |
| | Metals VOCs by 524.2, MCP 14 | 20 200 00 | | | | | | | | | | | | | | | analytes. The applicable MCP reporting limits were achieved. CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method |
| MW-3 | Metals | 11-Jun-10 | | Х | | Х | Х | | N | Y | Y | Υ | Υ | NA | | Υ | 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection lithan Method 8260. (Data is "non-CAM", not "CAM non-compliant.") |
| RIZ-3 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | х | | N | Υ | Υ | Υ | Υ | NA | | Υ | CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection lithan Method 8260. (Data is "non-CAM", not "CAM non-compliant.") |
| MW-2 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | х | | N | Υ | Υ | Υ | Υ | NA | | Υ | CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection lithan Method 8260. (Data is "non-CAM", not "CAM non-compliant.") |
| GHC-6 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | х | | N | Y | Υ | Y | Υ | NA | | Y | CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Methor 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection lithan Method 8260. (Data is "non-CAM", nor "CAM non-compliant.") |
| RIZ-9 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | Х | | N | Υ | Υ | Υ | Υ | NA | | Y | CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection lithan Method 8260. (Data is "non-CAM", not "CAM non-compliant.") |
| RIZ-10 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | х | | N | Y | Y | Y | Υ | NA | | Y | CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection li |
| RIZ-8 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | Х | | N | Y | Y | Y | Υ | NA | | Y | than Method 8260. (Data is "non-CAM", not "CAM non-compliant.") CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Method 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection li |
| MW-9 | VOCs by 524.2, MCP 14 Metals | 11-Jun-10 | | х | | х | х | | N | Y | Y | Y | Υ | NA | | Y | than Method 8260. (Data is "non-CAM", not "CAM non-compliant.") CAM Compliant is N because VOC analysis was performed by EPA Method 524 rather than EPA Method 8260. Methor 524 was used to be consistent with the requirements of the annual BOH sampling and because it has lower detection li |
| | Ivietals | | | | | | | | | | | | | | | | than Method 8260. (Data is "non-CAM", not "CAM non-compliant.") |
| RIZ-9 (9'-11') | VOCs by 8260 (High Range and Low Range), MCP 14 Metals | 6-Dec-07 | х | | | х | х | | Υ | Y | Υ | Υ | NA | NA | | Υ | LCS/LCSD % recoveries for Dichlorodifluorometh are below the individual acceptance critera for the compounds, but within overall method allowances. |
| RIZ-10 (44'-46') | VOCs by 8260 (High Range and Low Range), MCP 14 Metals | 6-Dec-07 | х | | | х | х | | Y | Y | Υ | Υ | NA | NA | | Υ | LCS/LCSD % recoveries for Dichlorodifluorometh are below the individual acceptance critera for the compounds, but within overall method allowances. |
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| Appendix B - Data Usability Ass | pendix B - Data Usability Assessment Summary Table | | | | | | | | | | | | | | | |
|---------------------------------|--|-----------------|----------|--------|----------|--------------------|-------|--------------|--------------|---------------------|---------------|-----------------|--------------------|-----------------------------|---------------------|---------------------|
| Sample ID or Series | Parameters | Date Sampled | 8 | Ground | Sedimen | Site Characteriza: | EPCS | CAM Compile. | Sample Reco. | Method QA/QC Proces | QA/QC Require | Trip Blanks OK? | Field Duplicate OK | Relative Percent Difference | Appropriate Reposit | Data Qualifications |
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| Appendix B - Data Usability As: | sessment Summary Table | | | | | | | | | | | | | | | |
|---------------------------------|------------------------|-----------------|------|------------|-----------------------|--------------|--------------|-----------|-----------------------|---------------|-----------------|---------------------|-----------------------|---------------|---------------|---------------------|
| Sample ID or Series | Parameters | Date Sampled | Soll | Groudwater | Sediment Site Char | Hazard Eliza | CAM Complian | Sample R. | Method QA/QC Procedur | QA/QC Require | Trip Blanks OK? | Field Duplicate OK? | Relative Percent Div. | Appropriate R | Limits? (Y/N) | Data Qualifications |
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Appendix C

Public Notification Documentation

from Norfolk to CNH.

Dec. 23 ansported one d Post Road to

Transported m Main Street

wo-car motor it reported on No injuries.

ec. 24 ansported one Iltop Drive to

ansported one oute 1 South d Hospital. ansported one ain Street via d Hospital.

Transported n Main Street vood Hospital.

Dec. 25

CO detector low battering Main Street.

Assisted the d out of the ain Street. ansported one ain Street via

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/est St.,

s by 12o, please

activation unintentional on Killeen Road. Smoke was from cooking. Alarms had resent prior to the Fire Department's arrival Investigation only.

10:01 p.m. Report of people being stuck in an elevator on Elm Street. They were selfrescued before the Fire Department arrived. Elevator removed from Building manager notified.

10:05 p.m. A small grease fire in the oven was reported on Endean Drive. The fire was out prior to the Fire Department's arrival. Smoke had to be ventilated from the home. After ventilation, the alarm was reset. No damage to the oven.

Sunday, Dec. 26

2:57 a.m. Responded to a report of a fire in the living

Norfolk a.m. Ambulance transported one person from Drake Circle via Domenica Road.

1:38 p.m. CO detector activation on Plain Street. No CO BLS = Basic Life Support CNH = Caritas Norwood

www.dedhamsavings.com

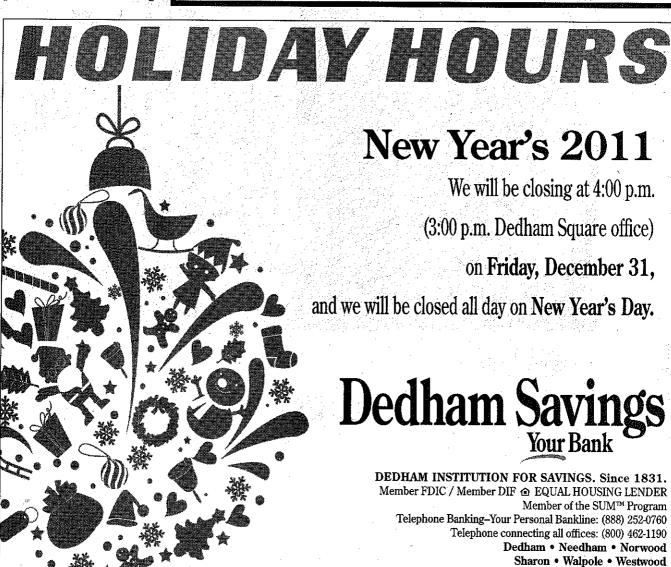
NOTICE OF A PUBLIC INVOLVEMENT PLAN MEETING

Walpole Park South Walpole Park South Road Walpole, Massachusetts RTN 3-21915

Walpole Park South received a petition from residents in Walpole requesting this location be designated as a Public Involvement Plan (PIP) site, in accordance with MGL c.21E § 14 (a). A PIP, dated April 6, 2005, was prepared for this Site indicating that meetings would be held to present major site documents to the public.

In accordance with the PIP, a public meeting will be held in the Main Meeting Room, Walpole Town Hall, 135 School Street, at 7:00 PM on January 13, 2011 to present the draft Phase V Completion Statement/Class B-1 Response Action Outcome Statement and to provide an opportunity for public comment. Copies of the draft report will be available at the meeting, and will also be placed in the public information repository at the Walpole Public Library on or before January 13, 2011. The public comment period for the RAO Statement will run through February 2, 2011.

Any questions regarding this meeting or the Public Involvement Plan should be directed to Raymond C. Johnson, P.G., L.S.P., Senior Vice President, Tetra Tech, Inc., 1 Grant Street, Framingham, MA 01701-9005, at 508-903-2000.



January 10, 2011

Mr. Christopher G. Timson, Chairman Board of Selectmen Town of Walpole 135 School Street Walpole, MA 02081

Re: Notice of Class A-2 Response Action Outcome Statement and Phase V Completion Report

Walpole Park South Walpole, Massachusetts

RTNs 4-3021915

Dear Mr. Timson:

On behalf of Walpole Park South, Tetra Tech, Inc. is providing this notification that a Class A-2 Response Action Outcome Statement and Phase V Completion Statement will be submitted to the Massachusetts Department of Environmental Protection (DEP) for the above referenced Disposal Site on or about February 3, 2011.

This notification is being made pursuant to the requirements of the Massachusetts Contingency Plan (MCP) 310 CMR 40.0000. The report will be available for review at the DEP Southeast Regional Office located at 20 Riverside Drive in Lakeville, Massachusetts by appointment. A copy of the report will also be available in the Public Information Repository at the Walpole Public Library.

Please contact the undersigned if you have any questions.

Very truly yours,

Raymond C. Johnson, P.G., L.S.P. Senior Vice President

P:\Pre-FY2008\12700000\12700058\12700058-003\RAO Statement\Town Notification Letter.doc

January 10, 2011

Ms. Robin Chapell, Health Agent Town of Walpole Board of Health 135 School Street Walpole, MA 02081

Re: Notice of Class A-2 Response Action Outcome Statement and Phase V

Completion Report Walpole Park South Walpole, Massachusetts RTNs 4-3021915

Dear Ms. Chapell:

On behalf of Walpole Park South, Tetra Tech, Inc. is providing this notification that a Class A-2 Response Action Outcome Statement and Phase V Completion Statement will be submitted to the Massachusetts Department of Environmental Protection (DEP) for the above referenced Disposal Site on or about February 3, 2011.

This notification is being made pursuant to the requirements of the Massachusetts Contingency Plan (MCP) 310 CMR 40.0000. The report will be available for review at the DEP Southeast Regional Office located at 20 Riverside Drive in Lakeville, Massachusetts by appointment. A copy of the report will also be available in the Public Information Repository at the Walpole Public Library.

Please contact the undersigned if you have any questions.

Very truly yours,

Raymond C. Johnson, P.G., L.S.P. Senior Vice President

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Appendix D **Soil Boring Logs and Monitoring Well Construction Diagrams**

| | Geo | Hyd | roCy(| CLE, I | NC. | | $\mathbf{E}_{\mathbf{i}}$ | nvironme | ntal Drill | ing Log |
|--------------|------------|-----------------------|----------------------------------|-------------|------|----------------------------------|---|---------------------------|------------------------------|--------------------|
| | 425 Newton | ville Avenue | Newton, MA 024 | 60 | | (617) 527-8074 (617) 527-8668 | | | | |
| | Project: | Walpol | e Park Sou | ıth | | | | Project No. GHC# 03027 | Location No. GHC-1 | Sheet 1 of 1 |
| | Drilling | Location | _ | _ | | first building toole Park Sout | to the west of the Route 1 h. | | 9/04 @ 14:50 9/04 @ 16:15 | |
| | Drill Ri | g: Truck | Mounted l | Rig | | Inspector: TV | /M and KAR | | | |
| | | | eter: 6.25" | | | 4 | - Steve and Tim | | ter Depth @ Co | 4 |
| | | Length: | plit Spoon | | | Temperature: | nny and windy 25° | Date/Time | Depth | Meas. Pt. TOC/Grnd |
| ď | | | Sample | | Rock | Soil | Sample Description/ | Stratigraphic | Mate | 1 |
| | Depth | No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | Insta | lled |
| | | | | | | | | | | |
| | | | · | | | | | | Road Box (0.0') | |
| | | SS-1 | (0-2') | 24" | | 46/41/38/31 | Dense, brown, f-c SAND, some Gravel, | | Cement 1-0' | |
| | | | | | | | Topsoil. | | | |
| | 5 | SS-2 | (5-7') | 12" | | 28/38/38/20 | Dense, brown, f. SAND, broken | | Betonite 3-2' | 7 |
| | | 55-2 | (3-7) | 12 | | 26/36/36/20 | rock pieces. | | 2" PVC Riser | |
| | | | | | | | | | Pipe 11.5-0' | |
| in the | 10 | SS-3 | (10-12') | 8" | | 12/33/41/17 | Dense, brown, f-c SAND, broken | | | |
| J | | | (10 10) | | | | rock pieces, Saturated. | | | |
| | | | | | | | | | Sand 20-3' | |
| eres i | 15 | SS-4 | (15-17') | 4" | . " | 6/10/12/15 | Medium, brown, f-c SAND, | | | |
| i | - | | | | | | broken rock pieces. | | 2" PVC 10-Slot | |
| es (d | | | | | | | | | Screen 20-5' | |
| | 20 | | | | | | 1 | | | |
| | | | | | | | EOB = 20 feet. | - | | 4 1 |
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| | | and some little | 35 to 50 20 to 35 10 to 20 | % % % | | 4-10 10-30 r | very loose Slight loose Low medium Medium | | LAY 1/8" SILT 1/16" | |
| | | trace | 1 to 10 | % | , | | dense High very dense Very H | Silty CLA ligh CLAY | Y 1/32" 1/64" | |

| J GEO | HYD | ro C yo | CLE,] | INC. | | E | nvironme | ntal Drill | ing Log |
|--------------|----------------|---------------------------------------|-------------|-------------|---|---|---------------------------------------|-----------------|--|
| 425 Newton | rville Avenue | Newton, MA 024 | 160 | | (617) 527-8074 (617) 527-8668 | | | | <i>U</i> - <i>O</i> |
| Project: | Walpol | le Park Sou | ıth | | | | Project No. | Location No. | Sheet |
| T | - · | | | 1 | | | GHC# 03027 | GHC-2 | 1 of 1 |
| Drilling | Locatio | | | | | e parking lot for the building | | 9/04 @ 13:00 | |
| Deili D: | or Two-1- | on the Mounted | e southw | estern s | | pole Park South Property. | Finished: 01/1 | 9/04 @ 14:40 | |
| | | eter: 6.25" | | | Inspector: TW Driller: TDS | | Canana david | ton Donth @ Co | 1_4: |
| 20 | | Split Spoon | | | | nny and windy | Date/Time | ter Depth @ Cer | Meas. Pt. |
| | Length: | | | | Temperature: | | 1/19/04 @ 15:35 | Depth 18.19 | TOC/Grnd |
| | T | Sample | | Rock | Soil | Sample Description/ | Stratigraphic | Mater | 1 |
| Depth | No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | Insta | |
| | | | | | | | | | 1 |
| | | | | | | | | | |
| | | | | | | | | Road Box | |
| | | | | | | | | (0.0') | |
| | | | | | | | | (0.0) | |
| | SS-1 | (0-2') | | | | Brown, f-m SAND, little Gravel. | | Cement 2-0' | |
| | | ` ′ | | | | | | 7 | |
| | | | | | | | | | |
| | | | | | | | | Betonite 5-2' | ₩ ₩ |
| 5 | SS-2 | (5-7') | 14" | | 12/38/48/60+ | Dense, brown, f-c SAND, trace Gravel. | | Botonic 3-2 | ₩ / ₩ |
| | | | | | | | | 2" PVC Riser | |
| | | | | | | | | Pipe 7-0' | |
| 4 ., | | | | | | | · | 1 ipe 7-0 | |
| | | | · | | | | | | |
| 10 | SS-3 | (10-12') | 0" | | 102/-/-/- | No Sample. | | | |
| | | (10 12) | | | 102/-/- | no sample. | | | · |
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| | | | | , | | | | G 100 51 | |
| | | | | , | | | · · · · · · · · · · · · · · · · · · · | Sand 22-5' | |
| 15 | SS-4 | (15-17') | 12" | | 29/19/30/15 | Medium, brown, f-m Sand, little Gravel, | | | - |
| | - 55 | (15 17) | - 12 | | 23/13/30/13 | Saturated. | | | - ∰≡=₩₩ |
| | | · · · · · · · · · · · · · · · · · · · | | | | Sauraicu. | | 28 DVC 10 GL-4 | |
| 7.5 | | | | | | | | 2" PVC 10-Slot | 1 = |
| | | | | | | | | Screen 22-7' | ₩ |
| 20 | SS-5 | (20-22') | 7" | | 55/80/-/- | Dense, brown, f-c SAND, trace Silt, | . , | | |
| | 55.5 | (20 22) | | | 33/60/ | | | | |
| 7 | | | | | | trace Gravel. | | | |
| | | | | | | FOR - 22 foot | | | |
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| and some | | io 30% io 35% | | | 4 very loose10 loose | Low | Clayey SILT SILT & CLA | 1/4" Y 1/8" | |
| little | | o 20% | | | -30 medium | Medium | CLAY & SIL | | |
| trace | 11 | to 10% | | 30- | | High | Silty CLAY | 1/32" | Ī |
| | | | | >: | 50 very dense | Very High | CLAY | 1/64" | |

| | Geo | Hyd | ROCYO | CLE,] | Inc. | | Eı | nvironme | ntal Drill | ing Log |
|---|----------|-----------------|----------------------|--------|----------|----------------------------------|--|---------------------------------|-----------------------|--|
| | | | Newton, MA 024 | | | (617) 527-8074 (617) 527-8668 | | | | Lot |
| | Project: | Walpol | e Park Sou | ıth | | | | Project No. GHC# 03027 | Location No. GHC-3 | Sheet 1 of 1 |
| | Drilling | Locatio | | | | | parking lot for the building | Begun: 01/1 | 9/04 @ 13:10 | 1 |
| | - '11 D' | m 1 | | | estern s | | pole Park South Property. | Finished: 01/1 | 9/04 @ 14:30 | |
| | | | Mounted leter: 6.25" | | · | Inspector: L Driller: TDS | | 1 Groundwe | ater Depth @ Co | mnletion |
| | | | Split Spoon | | | <u> </u> | nny and windy | Date/Time | Depth (a) Co | Meas. Pt. |
| | | Length | 24" | | | Temperature: | 25° | | | TOC/Grnd |
| | | | Sample | · . | Rock | Soil | Sample Description/ | Stratigraphic | Mate | |
| | Depth | No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | Insta | lled |
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| | | SS-1 | (0-2') | 24" | | 32/32/11/5 | Mediun, brown, f. SAND, some Gravel, | | Cement 1-0' | |
| | | | | | | | Topsoil. | | Madina Pill 0 41 | ************************************** |
| | | | | | | | | | Native Fill 2-1' | |
| | 5 | SS-2 | (5-7') | 12" | | 14/26/20/11 | Medium, brown, f. SAND, some Gravel. | | Betonite 4-2' | 7 |
| | | | | | | | | | 2" PVC Riser | |
| | | | | | | | | , | Pipe 6-0' | |
| | | - | | | | | | \ | | √ ■ · · |
| | 10 | SS-3 | (10.120) | 18" | | 12/12/19/15 | Madium Lucum & CANTO little Count | | Sand 16-4' | : = |
| | 10 | 33-3 | (10-12') | 18 | | 13/13/18/15 | Medium, brown, f. SAND, little Gravel. | | | · 📜 |
| | | | | | | | | | 2" PVC 10-Slot | 1 |
| | | | | | | | | | Screen 16-6' | 1 🗏 |
| | | | | | | · | | , | | |
| | 15 | SS-4 | (15-17') | 10" | | 13/100+/-/- | Dense, brown, f. SAND, | | | - |
| | | | | | | | broken rock pieces. EOB = 16 feet. | + | | |
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| | | | papara T | · • | | | | High CLAY | | |

| | GEO | HYD | roCy | CLE, | INC. | | F | nvironme | ental Drill | ino I | ഹ |
|------|----------------|-----------|------------------|----------------------|---------------------------------------|----------------------------------|--------------------------------------|--------------------|----------------|-------------|--------|
| | l | | Newton, MA 024 | | | (617) 527-8074 (617) 527-8668 | | | | mg L | νg |
| | Project: | Walpol | le Park Sou | uth | · · · · · · · · · · · · · · · · · · · | (017) 327-0008 | | Project No. | Location No. | Sheet | , |
| | Drilling | Location | n. I costo | din a fia | 1.1 : 41 | | | GHC# 03027 | GHC-4 | 1 of | 1 |
| | Bumin | Locatio | II. Located | o m a ne South Pi | oid in the | nortneastern i | nost portion of the Walpole | | 19/04 @ 11:20 | | |
| | Deill Ri | g: Truck | Mounted | Dia | roperty. | IIngmostow TV | TA A | Finished: 01/1 | 9/04 @ 12:50 | | |
| | Drill H | de Diam | eter: 6.25" | rug | | Inspector: TV | | T . | | | |
| | | | Split Spoor | | | Driller: TDS | | 1 | ater Depth @ 🗞 | 7 | |
| | | r Length: | | 1 | | weather: St | nny and windy | Date/Time | Depth | Meas. | |
| | Bampic | T Lengur. | | | T | Temperature: | | 1/19/04 @ 15:05 | 11.87 | TOC/G | rnd |
| | Danth | | Sample | | Rock | Soil | Sample Description/ | Stratigraphic | Mate | | |
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| ۷ | | SS-1 | (0-2') | | | | Light-brown, f. SAND, trace c. Sand. | | Cement 1-0' | | |
| | | | | | | | 1 | | | - 123 | e leva |
| | | | | | | | | | Betonite 2-1' | J00000 / | 00000 |
| | | | | | | | 1 | | Belomic 2-1 | <i> </i> | |
| | 5 | SS-2 | (5-7') | 16" | | 17/20/17/16 | Medium, brown, f-c SAND, little | 1 | | H/\square | |
| 3558 | | | | | | | Gravel. | 1 | OR DATE D | $H \equiv$ | |
| | | | | | | | Gravei. | | 2" PVC Riser | ∤ <u> </u> | |
| 1000 | | | | | | | , | | Pipe 3-0' | ∤ | |
| | | | | | | | | | | | |
| I | 10 | SS-3 | (10-12') | 14" | | 16/17/10/0 | . | | Sand 13-2' | | |
| | 10 | 33-3 | (10-12) | 14" | | 16/17/18/3 | Medium, dark-brown, f-m SAND, | | | | |
| | | | | | | | Saturated. | | | | |
| | | | | | | | | | 2" PVC 10-Slot | | |
| | | | | | | | | | Screen 13-3' | | |
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| | some little | | to 35% | | | 4-10 loose 10-30 mediur | Low | SILT & CLAY | 1/8" | | |
| | trace | | to 20% to 10% | | | 10-30 mediur 30-50 dense | n Medium High | CLAY & SILT | 1/16" | | 1 |
| | 440 | | 10/0 | | 1. | >50 very de | | Silty CLAY CLAY | 1/32" 1/64" | | - 1 |
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| $\mathbf{J}\mathbf{G}_{\mathrm{E}}$ | OHYD | ROCY | CLE, | INC. | | E | nvironme | ntal Drill | ing Log |
|-------------------------------------|------------|---------------|-----------|-----------|----------------|--|------------------|--------------------|---------------------|
| 7 | | Newton, MA 02 | | | (617) 527-8074 | | | | 00 |
| Proje | ct: Walpo | le Park So | uth | | (617) 527-8668 | | Project No. | Location No. | Sheet |
| 7 | | | | | | | GHC# 03027 | GHC-5 | 1 of 1 |
| Drilli | ng Locatio | n: Locate | d in nort | hwestern | portion of the | e parking lot for the building | | 9/04 @ 09:00 | |
| 15.211 | D: 70 1 | on th | e southy | vestern s | | pole Park South Property. | Finished: 01/1 | 9/04 @ 11:10 | |
| Drill . | Kig: Truck | Mounted | Rig | | Inspector: T\ | | · | | |
| | | neter: 6.25" | | | Driller: TDS | | Groundwa | ter Depth @ Co | mpletion |
| | | Split Spoor | 1 | | | unny and windy | Date/Time | Depth | Meas. Pt. |
| Samp | ler Length | | | Ι | Temperature | · | | | TOC/Grnd |
| n | | Sample | 1 | Rock | Soil | Sample Description/ | Stratigraphic | Mate | |
| Dept | h No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | Insta | lled |
| | | | | | | | | Road Box (0.0') | € |
| | SS-1 | (0-2') | | | | Brown, f-c SAND, trace Gravel, trace | | Cement 2-0' | |
| | | | | | | Silt. | | Comone 2-0 | |
| 7 | | | | | | | | | |
| | | | | | | 1 | | Native Fill 6-2' | / / / |
| 5 | SS-2 | (5-7') | 9" | | 9/7/6/16 | Very loose, Grey, f-c SAND, trace | | Nauve Fill 0-2 | } |
| | | () | | | 3,7,0,10 | • | | A | / |
| | | | | | | Gravel, trace Silt. | | 2" PVC Riser | 6000 00000 |
| | - | | | | | · · | | Pipe 10-0' | |
| | | | - | | | 4 | | | 18888 188888 |
| | | | | | | | | Betonite 8-6' | |
| 10 | SS-3 | (10-12') | 17" | | 11/5/9/14 | Very loose, Grey-brown, f SAND, trace | | | |
| | | | | | | m Sand, trace c Sand, trace Silt. | | | |
| 7 | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | 1 | | | 18 🖃 📟 |
| 15 | SS-4 | (15-17') | 13" | | 7/14/14/21 | Loose, Brown, f-c SAND, trace Gravel, | | · | |
| | | | | | | trace Silt, Saturated. | | | |
| - | | | | | | Samuelou. | | | |
| | | | | | | | , · | G140.0t | / 🗐 📗 |
| 7 | | | | | * | | | Sand 40-8' | |
| 20 | SS-5 | (20-22') | 17" | | 7/12/21/10 | ACT D CONT | | | |
| 20 | 33-3 | (20-22) | 17 | | 7/12/21/18 | Medium, Brown, f-c SAND, trace | | | |
| , | · | | | | | Gravel, trace Silt, Saturated. | · | | |
| | | | | | | | | | |
| | | | | | | 1 | | | |
| | - | | | | | 1 | | | |
| 25 | SS-6 | (25-27') | 14" | | 6/19/31/34 | Medium, Grey, f-c SAND, trace Silt, | | | |
| | | | | | | Saturated. | | | |
| | |] | | | • | | | | |
| ľ | | | | | | | | 2" PVC 10-Slot | ľ⊟ I |
| | | , | | | | | | Screen 40-10' | |
| 30 | SS-7 | (30-32') | | | 25/32/46/38 | Dense, Grey-brown, f SAND, trace | | | |
| | | | | . 1 | | m Sand, trace c Sand, trace Silt, | | | |
| | | | | | | Saturated. | | | |
| | | | | | | | | | |
| | | | | | | Dense, Grey-brown, f SAND, trace | | | |
| 35 | SS-8 | (35-37') | 17" | | 24/38/46/66 | m Sand, trace c Sand, trace Silt, Moist. | | | |
| | | OMPONENT | | | | | ATT DI ACTUCIONE | | |
| | and | 35 to 509 | | | 0-4 | very loose Sligh | t Clayey | SILT 1/4" | |
| | some | 20 to 359 | | | 4-10 | loose Low | | CLAY 1/8" | , |
| 1 | little | 10 to 20% | | | 10-30 | medium Medi | | | |
| | trace | 1 to 109 | % | | 30-50 | dense High | Silty Cl | | |
| | | | | | >50 | very dense Very | High CLAY | 1/64" | İ |

| - | | | | | (617) 527-8668 | | | | |
|----------|-----------------|-----------|--------|----------------|---|--|-------------------|----------------|--|
| roject: | Walpol | ļ. | | | | | Project No. | Boring No. | Sheet |
| | | | | , | *************************************** | | GHC# 03027 | GHC-5 | 2 of |
| | | Sample | | Rock | Soil | | Stratigraphic | Mate | rials |
| epth | No. | Depth | Recov. | RQD | Blows per 6" | Sample Description | Description | Insta | alled |
| | | | | | | · | | . 6 | |
| - 1 | | | | | | | | 2" PVC 10-Slot | |
| | | | | | | Dense, Brown, f SAND, little Silt, trace | | Screen 40-10' | |
| | | (10.10) | | - | | Gravel, trace m Sand, trace c Sand, | | | |
| 40 | SS-9 | (40-42') | 12" | | 28/48/29/19 | Moist. | | | |
| ŀ | | | | | | EOB = 40 feet. | | | - |
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| 80 | | | | | | | est of the second | |]_] |
| | | OMPONENT | | | | RELATIVE DENSITY | OVERALL PLAST | ICITY | |
| | and | 35 to 50% | | | • | 0-4 very loose | Slight | Clayey SILT 1/ | 4" |
| | some | 20 to 35% | | | | 4-10 loose | | | /8" |
| | little trace | 10 to 20% | | | | 10-30 medium 30-50 dense | | | 16" |
| | | | | | | | High | Silty CLAY 1/3 | |

| | GEO | Hyd | roCyc | CLE, | INC. | | E | invironme | ntal Drill | ling Log |
|------|----------|----------|----------------|---------|--|-------------------|--|---------------------------------------|------------------|--------------------------------|
| 4 | | | Newton, MA 024 | | | (617) 527-8074 | | | | |
| | Project: | Walpo | le Park Sou | ıth | | (617) 527-8668 | | Project No. | Location No. | Sheet |
| 4 | | | | | | | | GHC# 03027 | GHC-6 | 1 of 1 |
| | Drilling | Location | | | | cnic tables in tl | ne center of the Walpole Park | | 9/04 @ 11:15 | |
| | | | | Proper | ty. | | | Finished: 01/1 | 9/04 @ 12:50 | • |
| | Drill Ri | g: Truck | Mounted 1 | Rig | | Inspector: LC | В | | | |
| T | Drill Ho | ole Dian | neter: 6.25" | | | Driller: TDS - | | Groundwa | ater Depth @ Co | mpletion |
| | Sampler | Type: S | Split Spoon | l | | Weather: Sur | | Date/Time | Depth | Meas. Pt. |
| | | Length | | | | Temperature: | | 1/19/04 @ 15:20 | 8.71 | TOC/Grnd |
| | | T S | Sample | | Rock | Soil | Sample Description/ | Stratigraphic | | erials |
| ı | Depth | No. | 1 | Γ. | 1 . | | Detector Readings | Description | 1 | alled |
| | Depth | No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | THE | alleu |
| | | | | | | | | | | |
| | | 1 | 1 | | 1 | | | | ļ | |
| | | | | | | 1 | | | Road Box | 1 .1 |
| | | | | | l | | | | (0.0') | 1 |
| | | | | | | | | | () | 1 |
| ŀ | | 66.1 | (0.00) | 04" | | 10/15/04/00 | \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ | | 0 10 | |
| | | SS-1 | (0-2') | 24" | - | 19/15/24/39 | Medium, brown, f-m SAND, | | Cement 1-0' | |
| | | | | | | ļ | some Gravel, Topsoil. | | i | 424 444 |
| T | | | | | | | | | Native Fill 2-1' | |
| | | L | | | | | | | Betonite 3-2' | Y Y [] |
| | 5 | SS-2 | (5-7') | 20" | | 9/23/26/34 | Medium, brown, f-c SAND, some | | | 7/= |
| | | | | | | | Gravel, broken rock pieces. | | 2" PVC Riser | -/ = |
| П | | | | | | | laraver, broken rock pieces. | | | 1 = 1 |
| | | | | | | | | | Pipe 4-0' | |
| | | | | | | | | | | 」/ 目 |
| | | | | | 1 | | | , | Sand 19-3' | |
| | 10 | SS-3 | (10-12') | 24" | | 7/8/12/14 | Loose, brown, f-c SAND, little gravel. | | | |
| | | | | | | | Saturated. | | | |
| | | | | | | | | | 2" PVC 10-Slot | |
| Ŧ | | | | <u></u> | | <u> </u> | | | | - = |
| | | | | | | | | | Screen 19-4' | - - - |
| | | | | | | | | | | |
| | 15 | SS-4 | (15-17') | 4" | | 60+/-/-/- | Dense, brown, f-m SAND, broken | | | |
| | | | | | İ | | rock, Saturated. | | | |
| | - | | | | | | | | | |
| | | | | | | | | | | |
| 7 | | | | | | | | | | |
| | 20 | | | | | | TOP 10.5 / | | | |
| | 20 | | | | | | EOB = 19 feet. | | | - |
| | | | | | | | | | | _ |
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| L | | | | | | | | | |] |
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| - FT | 25 | | | | | | | 1 | | - |
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| 4 | ł | | | - | - | | | | | - |
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| 7 | Į. | , | | | | | | | | _ |
| L | 35 | | | | | | | 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 | | |
| | 1 | MINOR C | OMPONENT | rs | | RELATI | VE DENSITY OVERA | ALL PLASTICITY | | |
| | | and | 35 to 50 | | | 0-4 | very loose Slight | | LT 1/4" | |
| T | | some | 20 to 359 | | | 4-10 | loose Low | SILT & C | | |
| | | little | 10 to 209 | | | 10-30 | medium Mediu | | | · |
| | | trace | 1 to 10 | % | | 30-50 | dense High | Silty CLA | | |
| | | | | | | >50 | very dense Very I | High CLAY | 1/64" | |

| | Geo] | Hydi | ROCYC | CLE, I | NC. | | E | nvironme | ntal Drilli | ng Log |
|-----|--------------------|----------|--|--|--|--|---|---------------------------------|--------------------------------------|--|
| | | | Newton, MA 0246 | • | | (617) 527-8074 (617) 527-8668 | | | | |
| | Project: | Walpole | Park Sou | th | | | | GHC# 03027 | GHC-7 | Sheet 1 of 1 |
| | Drilling | Location | | in a gra South Pr | | in the southea | stern portion of the Walpole | Begun: 01/19 Finished: 01/19 | 9/04 @ 09:1 0 9/04 @ 10:45 | |
| | | | Mounted I | | | Inspector: LC | | | | |
| 1 | | | eter: 6.25" | | | Driller: TDS - | | | ter Depth @ Co | Meas. Pt. |
| | Sampler Sampler | Type: S | plit Spoon | | | Temperature: | nny and windy | Date/Time 1/19/04 @ 15:13 | Depth 16.81 | TOC/Grnd |
| 4 | Sampler | Lengui. | Sample | | Rock | Soil | Sample Description/ | Stratigraphic | Mater | |
| | Depth | No. | Depth | Recov. | RQD | Blows per 6" | Detector Readings | Description | Insta | lled |
| No. | | | | | | | | | | |
| 4 | | | | | | | | | | ŧ |
| | | | | | | | | | Road Box | |
| | | | | | | | | | (0.0') | Λ |
| ٦ | | | | | | • | | | · | |
| | | SS-1 | (0-2') | 15" | | 143/172/-/- | Very dense, brown, f-m SAND, trace | | Cement 1-0' | |
| | | | | | | | Gravel. | | | 4 1 1 1 |
| 7 | | | | | | | | | | 4 |
| | | | | | | | | | 27.1. 77.11.6.11 | 4/ |
| | 5 | SS-2 | (5-7') | 15" | | 8/9/10/10 | Loose, brown, f. SAND. | | Native Fill 6-1' | 4 1 4 1 |
| | | | | | | <u> </u> | | | | 80000 |
| | | | | | | <u> </u> | | | | -‱ ‱ |
| | | | | | <u> </u> | | 4 | | D-4i4- 9 61 | _ |
| | | | | | - | | l | | Betonite 8-6' | 1 / |
| | 10 | SS-3 | (10-12') | 16" | | 7/9/12/10 | Medium, brown, f. SAND. | | | - /= |
| | | | | ļ | <u> </u> | | | | 2" PVC Riser | -/ 目 □ |
| | | | | | <u> </u> | | 4 | | Pipe 10-0' | ┪ |
| | | <u> </u> | - | | | - | - | | 1 ipe 10-0 | - |
| | 15 | SS-4 | (15-17') | 16" | | 3/6/3/3 | Loose, brown, f. SAND, Saturated. | | | 1 🗐 🛚 |
| | 1.13 | 33-4 | (13-17) | 10 | 1 | 310/3/3 | Loose, blown, i. SAND, Saturated. | | | - |
| | | | | | | | <u> </u> | | | 1,= |
| | | | | | | 7 | | | Sand 25-8' | 7 🗐 🛮 |
| | - | | | | <u> </u> | | | | | |
| | 20 | SS-5 | (20-22') | 24" | | 10/10/31/16 | Medium, brown, f-m SAND. | | : | |
| | , | | | | , | | | • | · | |
| ٦ | l | | | | | | | | 2" PVC 10-Slot | |
| | 1 | | | | | | | | Screen 25-10' | |
| | | | | | | | | | | _ |
| | 25 | SS-6 | (25-27') | 24" | | 18/30/12/13 | Medium, brown, f-c SAND, some | | | |
| | 1 | | | | | | Gravel. | | | _ |
| | | | | | | | EOB = 25 feet. | | | _ |
| | | | ļ | | _ | | | | | \dashv \mid \mid |
| | 1 | | | <u> </u> | <u> </u> | | | | | - |
| | 30 | | | | | | | | | - |
| | | | <u> </u> | <u> </u> | _ | - | 4 | • | | - |
| | 1 | | | <u> </u> | | . | ┥ | | | - |
| | | - | | | 1 | | - | | | \dashv $\mid \cdot \mid$ |
| | 35 | | | | 1 | | 1 | | - | - |
| | 33 | MINOP (| COMPONEN | JTS | 1 | אַ אַ אַ אַ אַ אַ אַ אַ אַ אַ אַ אַ אַ א | I VE DENSITY OVE | RALL PLASTICITY | | |
| | | and | 35 to 5 | | | 0-4 | very loose Sligh | ht Clayey S | | |
| | , | some | 20 to 3 | 5% | | 4-10 | loose Low | | | |
| | | little | 10 to 2 | | | 10-30 30-50 | medium Med dense High | | | |
| | | trace | 1 10 1 | V /0 | | >50 >50 | | High CLAY | 1/64" | |
| - | • | | | | | - | - · · · · · · · · · · · · · · · · · · · | | | |

| | R I Z Z A S S A TECH C | 0 C I | ATES | | | W | ELL | NUMBER RIZ-1 PAGE 1 OF 1 |
|------------------|-------------------------------------|-----------------|---|----------------|--|---|-----------|---|
| CLIEN | IT Walp | ole Pa | | | | | | |
| | ECT NUM | | | | | PROJECT LOCATION Walpole, Massac | | |
| | | | | | | GROUND ELEVATION GROUND WATER LEVELS: | HOLES | 51 2 E <u>2"</u> |
| | | | | | er | | | |
| LOGG | ED BY | Dimitri | Gounis | | CHECKED BY | AT END OF DRILLING | | |
| NOTE | s | 1 | ı | | I | AFTER DRILLING | | 1 |
| O DEPTH (ft) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | MA | ATERIAL DESCRIPTION | PID (ppm) | WELL DIAGRAM |
| | | 50 25 100 | 11-15-52- 93 (67) 42-30 11-42-40- 45 (82) | | Tan medium to fine sand, 12.0 12.5 boulder Tan fine sand with some part of the sand with some part o | plasticity with gravel and very course sand, wet, | 0.8 | Backfill: Cuttings PVC Riser Bentonite Filter Pack PVC Screen |
| - 25 - 30 | | | | | Be | ottom of hole at 32.0 feet. | | |

| | RIZZ(ASS) RATECH C | 0 C I | ATES | | | V | /ELL | NUMBER RIZ-2 PAGE 1 OF 1 |
|---|--------------------------------|---------------------------------|-----------------------------|------------------|---|--|-----------|---|
| | NT Walpo | | rk South Tru | | | PROJECT NAME Walpole Park South PROJECT LOCATION Walpole, Massac | husetts | |
| DATE DRILL DRILL LOGG | STARTE LING CON LING MET | D 10 ITRAC HOD Dimitri | TOR Soil Hollow Ste | Explora m Aug | | GROUND ELEVATION GROUND WATER LEVELS: AT TIME OF DRILLING | HOLE S | |
| o DEPTH | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | MAT | TERIAL DESCRIPTION | PID (ppm) | WELL DIAGRAM |
| | | 33 | 50 | | Tan madium to fine sand ar | nd gravel, dry, no odor | 0 | Backfill: Cuttings Bentonite PVC Riser |
| - 5 | | 0 | | ο · · · · | Gravel and cobbles off augo fine dry sand 7.0 | er, some angular gravel, some tan medium to | | |
| 10 | | 33 | 6-12-22-12 (34) | | Tan coarse sand and grave | l with fine sand, wet, no odor | 0 | Filter Pack PVC Screen |
| _ 15 _ | | 25 | 70 | | Tan coarse sand and grave odor, auger refusal at 15.5' 17.0 | el with fines, slightly plastic/cohesive, wet, no | 0 | |
| GENERAL BH/TP/WELL BORING LOGS.GPJ GINT US.GDT 3/15/06 2 2 2 5 6 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 | | | | | Bot | tom of hole at 27.0 feet. | | |

WELL NUMBER RIZ-3 RIZZO PAGE 1 OF 2 ASSOCIATES A TETRA TECH COMPANY CLIENT Walpole Park South Trust PROJECT NAME Walpole Park South PROJECT LOCATION Walpole, Massachusetts PROJECT NUMBER DATE STARTED 10/14/05 COMPLETED 10/14/05 GROUND ELEVATION HOLE SIZE 2" GROUND WATER LEVELS: DRILLING CONTRACTOR Soil Exploration DRILLING METHOD Hollow Stem Auger AT TIME OF DRILLING _---LOGGED BY Dimitri Gounis CHECKED BY AT END OF DRILLING _---**NOTES** AFTER DRILLING _---SAMPLE TYPE NUMBER BLOW COUNTS (N VALUE) GRAPHIC LOG (mdd) RECOVERY DEPTH (ft) MATERIAL DESCRIPTION WELL DIAGRAM Ы 0 Brown uniform medium sand fill, moist, no odor, off auger 3.0 5 5.0 Brown uniform medium to fine sand, moist, no odor 5-7-8-6 (15)Backfill: Cuttings 10 **PVC Riser** Brown/tan coarse sand and fine sand with gravel, wet, no odor 6-7-9-12 50 0 (16)15 Brown/tan coarse sand and fine sand with some gravel, wet, no odor 1-6-11-12 33 0 (17)Bentonite 20 Brown/tan coarse sand with some fines, wet, no odor 3-6-8-21 0 (14)**PVC Screen** 25 Filter Pack 15-20-23-Brown/tan coarse sand to very corase sand and gravel with some fines, 27 wet, no odor (43)Brown/tan coarse sand to very corase sand and gravel with fines, wet, no 7-11-9-11 83 0 (20)

GENERAL BH / TP / WELL BORING LOGS.GPJ GINT US.GDT 3/15/06

WELL NUMBER RIZ-3

RIZZO

GENERAL BH / TP / WELL BORING LOGS.GPJ GINT US.GDT 3/15/06

| PROJECT NUMBER PROJECT LOCATION Walpole, Massachusetts PROJECT LOCATION Walpole, Massachusetts A | ASSOCIATES A TETRA TECH COMPANY | | | PAGE 2 OF 2 | | | | |
|--|---------------------------------|------|---|--------------|--|--|--|--|
| HEDGO WELL DIAGRAM MATERIAL DESCRIPTION (a) Wounding prevents acurate sample from this depth, no soil desc. End or boring at 40', no refusal. | | | | | | | | |
| | | | | WELL DIAGRAM | | | | |
| | | 42.0 | 0 | | | | | |

| | RIZZ(ASS) | 0 C I | ATES | | | BORING NUMBER RIA PAGE 1 | |
|--|---|------------|-----------------------------|----------------|---|--|-----------|
| CLIEN | NT Walpo | ole Pa | rk South Tru | ust | | PROJECT NAME Walpole Park South | |
| | ECT NUM | | | | COMPLETED 10/14/05 | PROJECT LOCATION Walpole, Massachusetts GROUND ELEVATION HOLE SIZE 2" | |
| | | | | | ation | | |
| DRILL | DRILLING METHOD Hollow Stem Auger AT TIME OF DRILLING | | | | | | |
| | GGED BY Dimitri Gounis CHECKED BY AT END OF DRILLING TES AFTER DRILLING | | | | | | |
| NOTE | | | | | | AFTER DRILLING | |
| O DEPTH (ft) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | | MATERIAL DESCRIPTION | PID (ppm) |
| | | | | | Brown medium to fine sand | d and gravel, dry, no odor, off auger | 0 |
| 5 | | 75 | 25-30-27- 23 (57) | | Brown medium to fine sand | d and gravel, dry, no odor | 0 |
| | | | | | | | |
| | | 33 | 17-27-27- 29 (54) | | Brown medium to fine sand Auger refusal at 13' 12.0 | d and gravel with some coarse sand few orange mottles, moist, no odor. | 0 |
| GENERAL BH / TP / WELL BORING LOGS.GPJ. GINT US.GDT 3/15/06 12 12 1 | | | | | | Bottom of hole at 27.0 feet. | |

| | RIZZ ASS RATECH C | 0 C I | ATES | | | BORING NUMBER RI | |
|---|-------------------------|------------|-----------------------------|----------------|--------------------------|---|-----------|
| PROJ | ECT NUM | IBER | | | | | |
| | | | | | | GROUND ELEVATION HOLE SIZE _6" | |
| | | | | | er | GROUND WATER LEVELS: AT TIME OF DRILLING | |
| | | | | | CHECKED BY | | |
| NOTE | s | • | • | | | AFTER DRILLING | |
| O DEPTH (ft) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | | MATERIAL DESCRIPTION | PID (ppm) |
| | | 50 | 8-9-12-9 (21) | l**** | Brown medium to fine uni | form sand, moist, no odor | 0 |
| 5 | | 25 | 16-18-20- | | Brown medium to fine uni | form sand with angular gravel, dry, no odor | |
| - | | 25 | 25 (38) | | 7.0 | | 0 |
| 10 | | 0 | 6-21-45-30 (66) | | No recovery, brown mediu | um to fine uniform sand with angular gravel off auger | 0 |
| | | | | | | | |
| | | 83 | 6-9-11-7 (20) | | Brown/tan medium sand u | uniform with little gravel, moist, no odor | 0 |
| 20 | | | | | | Bottom of hole at 20.0 feet. | |
| ١٥٥.٥٥ الا ١٥٠.٥٥ الله | | | | | | | |
| LOGO.GFJ GI | | | | | | | |
| WELL BURING | | | | | | | |
| GENERAL BH / 1P / WELL BORING LOGS.GPJ GINI US.GD 3/19/08 | | | | | | | |
| S S S | | | | | | | |

| | RIZZ ASS | 0 C I | ATES | | | BORING NUMBER RIZ | |
|---|-----------------------|------------|-----------------------------|----------------|---------------------------------|--|-----------|
| PROJ | IECT NUI | /IBER | | | | | |
| DRILI | LING CO | NTRAC | TOR Geo | search | COMPLETED <u>2/16/06</u> | GROUND ELEVATION HOLE SIZE 6" GROUND WATER LEVELS: AT TIME OF DRILLING | |
| | GED BY | | | | CHECKED BY | AT END OF DRILLING AFTER DRILLING | |
| O DEPTH | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | | MATERIAL DESCRIPTION | PID (ppm) |
| | | 100 | 6-10-9-10 (19) | | 12" loamy top soil, 12" brown | medium to coarse sand and gravel with some fines, moist to wet, no odor. | 0 |
| <u>5</u> | | 33 | 10-9-8-7 (17) | | Brown medium to fine sand a | nd gravel with some organic material (plant), wet at bottom of spoon | 2.9 |
| 10 | | 83 | 3-9-10-13 | | Tan uniform medium to fine s | and, moist, no odor, boring ended due to proximity to overhead utilities, no | 0 |
| - | | | (19) | | 12.0 | Bottom of hole at 12.0 feet. | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| GENERAL DITTEN WELL BONING ECOS. O'S GINT GO. GDT STAND | | | | | | | |
| פבואבאטר בי | | | | | | | |

| | RIZZ (ASS (| 0 C I | ATES | | | BORING NUMBER RIA | |
|--|------------------------|------------|-----------------------------|----------------|---|---|-----------|
| PROJ DATE DRILI | ESTARTE | IBER O | 16/06 CTOR <u>Geo</u> | search | COMPLETED <u>2/16/06</u> | GROUND ELEVATION HOLE SIZE 6" GROUND WATER LEVELS: | |
| LOGO | | Chris I | Nitchie | | CHECKED BY | | |
| O DEPTH (ft) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | GRAPHIC LOG | | MATERIAL DESCRIPTION | PID (ppm) |
| | | | | | Tan to light brown medi | ium sand and gravel with some fines, wet (snow melt), no odor | |
| _ 5 | | 25 | 5-5-6-7 (11) | | Brown medium to coars | se sand with fines and some gravel, wet (snow melt), no odor | 0 |
| 10 | | 42 | 14-43-63- 30 (106) | | Tan/Brown medium to o | caorse sand with fines and gravel, wet, no odor. Likely near top of water table | 0 |
| 15 | - | | (106) | <u> </u> | 12.0 | | |
| | | 50 | 51 | | Tan/Brown uniform coa refusal at 18' 17.0 | rse sand with some fines transitioning to angular gravel with fines. Auguer | 0 |
| GENERAL BH / TP / WELL BORING LOGS.GPJ GINT US.GDT 3/15/06 | | | | | 17.0 | Bottom of hole at 18.0 feet. | |

WELL NUMBER RIZ-8 PAGE 1 OF 1



| PROJ | ECT NUM | BER . | rk South Tru 12700058 //5/07 | | | PROJECT LOCATION Walpole, Massachusetts | 2" |
|--------------|-----------------------|------------|------------------------------------|---------------------------------------|----------------|--|-------------------------|
| | | | TOR Geos | | | GROUND WATER LEVELS: | |
| | | | Hollow Ster | | | AT TIME OF DRILLING 24' | |
| | | | | CHECKED BY | | | |
| | S | | шоу | | | AFTER DRILLING | |
| IVOIL | | | <u> </u> | | | ALLENDRILLING | |
| O DEPTH (ft) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | REMARKS | GRAPHIC LOG | MATERIAL DESCRIPTION (교 d) 이 다 | VELL DIAGRAM |
| | S-1 | 50 | 6-6-8-5 (14) | | <u> </u> | gravei | Flush Mounted Road |
| - | S-2 | 90 | 5-3-4-7 (7) | | | .0 | Box |
| _ | | | | | | Dry, light tan, well sorted fine sand | |
| 10 | S-3 | 80 | 6-13-21-18 | Sample taken and submitted for VOC | | 0.0 Dry, light tan, well sorted fine sand | |
| - | | | (34) | and MCP-14 analysis | | Dry, gray, large gravel, little sand | |
| - | S-4 | 80 | 12-23-23- 30 (46) | from 9'-11' (RIZ-8-9'-11') | | 4.0 6.0 Dry, Brown, poorly sorted coarse sand, some silt, rocky | ✓2" PVC Capped Riser |
| 20 | | | 11-18-15- | | •.•ত্যু• | 9.0 Dry, tan, poorly sorted coarse and fine sand, some 0.1 | |
| | S-5 | 80 | 17 | | <u></u> | Dry, tan, poorly sorted coarse and fine sand, some 0.1 gravel, rocky | |
| _ | | | (33) | | | 4.0 | |
| | S-6 | 50 | 14-23-27- | | | Wet, brown, large gravel and rocks with coarse 0.1 | |
| | ' | | 21 (50) | | • • • • • | sand. Auger refusal at 27', bedrock | |
| 30 | | | | | | | ■Bentonite Seal |
| _ | | | | | | [73] | 8.3 |
| _ | | | | | | | |
| _ | | | | | | | |
| _ | | | | | | | |
| 40 | | | | | | | |
| _ | | | | | | | |
| _ | | | | | | | |
| - | | | | | | | |
| - | | | | | | | |
| 50_ | | | | | | | |
| - | | | | | | | |
| - | | | | | | | Sand Filter |
| - | | | | | | | Pack |
| | | | | | | | |
| 60 | | | | | | | |
| - | | | | | | | |
| - | | | | | | | 割 |
| - | | | | | | | |
| | | | | | | | 13 |
| 70_ | | | | | | | |
| - | | | | | | | ✓2" Machine |
| | | | | | | | Slotted Well Screen |
| _ | | | | | | | 2" PVC Plug |

| Tt TI | ETRA TECH R | IZZO | V | VELL NUMBER RIZ-8S PAGE 1 OF 1 |
|---------------------------------|----------------------|------------|---------------------------------|---|
| CLIENT Wa | alpole Park South 1 | Trust | PROJECT NAME Walpole Park South | 1 |
| PROJECT N | JMBER <u>1270005</u> | 8 | PROJECT LOCATION Walpole, Mass | sachusetts |
| | | | GROUND ELEVATION | HOLE SIZE 2" |
| DRILLING C | ONTRACTOR Ge | eosearch | GROUND WATER LEVELS: | |
| DRILLING M | ETHOD Hollow S | item Auger | AT TIME OF DRILLING N/A | |
| LOGGED BY | Luke Tulley | CHECKED BY | AT END OF DRILLING | |
| NOTES | | | AFTER DRILLING | |
| DEPTH (ft) SAMPLE TYPE NI IMBER | GRAPHIC | MATE | ERIAL DESCRIPTION | WELL DIAGRAM |
| 0 | | | | Flush Mounted Road Box |
| 5 - | | | | |
| 10 _ | | | | ■ Bentonite Seal |
| 15 | | | | Sand Filter Pack 2" Machine Slotted Well Screen |

WELL NUMBER RIZ-9 PAGE 1 OF 1

| CLIEN | IT Walp | ole Pa | rk South Tru | st | | PROJECT NAME Walpole Park South | | | | |
|---|-----------------------|-------------|-----------------------------|--|--|---|------------------------------|--------------------------------------|--|--|
| PROJ | ECT NUM | MBER . | 12700058 | | | PROJECT LOCATION Walpole, Massachu | ATION Walpole, Massachusetts | | | |
| DATE | STARTE | D 12 | 2/6/07 | COMPLETED | 12/6/0 | 7 GROUND ELEVATION | HOLE S | SIZE _2" | | |
| DRILL | ING CON | NTRAC | TOR Geos | search | | GROUND WATER LEVELS: | | | | |
| DRILL | ING MET | HOD | Hollow Ster | m Auger | | AT TIME OF DRILLING 14' | | | | |
| LOGG | ED BY _ | Luke 1 | Γulley | CHECKED BY | | AT END OF DRILLING | | | | |
| NOTE | s | | | | | AFTER DRILLING | | | | |
| о ДЕРТН (#) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | REMARKS | GRAPHIC LOG | MATERIAL DESCRIPTION | PID (ppm) | WELL DIAGRAM | | |
| - | S-1 | 50 | 3-2-3-4 (5) | | <u> </u> | gravel | 0 | Flush Mounted Road Box | | |
| 5 - | S-2 | 15 | 9-10-12-18 | | | 4.0 | 0 | | | |
| | 5-2 | 15 | (22) | | \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ | 9.0 | 0 | | | |
| 10 | S-3 | 80 | 3-5-9-10 (14) | Sample taken and submitted for VOC and MCP-14 analysis from 9'-11' | | Dry, tan uniform coarse sand with some fines transitioning to angular gravel with fines | 0 | Capped Riser | | |
| | | | | (RIZ-9-9'-11') | •••••• | 14.0 Wet, tan, well sorted coarse sand with angular | | - | | |
| _ 15 _ | S-4 | 50 | 34-19-13-7 (32) | | | gravel. Soft bedrock at 16'. | 0 | - | | |
| 20 | | | | | | | | ■ Bentonite Seal | | |
| 0 LOGS.GPJ 0 | | | | | | | | Sand Filter Pack | | |
| GENERAL BH / TP / WELL BORING LOGS. GP.J. GINT US.GDT 12/11/07 S | | | | | | | | 2" Machine Slotted Well Screen | | |
| GENERAL 35 | | | | | | Bottom of hole at 35.0 feet. | | 2" PVC Plug | | |

WELL NUMBER RIZ-10 PAGE 1 OF 1

| TETRA TECH RIZZO |
|------------------|
|------------------|

| CLIENT _Walpole Park South Trust PROJECT NUMBER _12700058 DATE STARTED _12/6/07 | | | | | | PROJECT NAME Walpole Park South | PROJECT NAME Walpole Park South PROJECT LOCATION Walpole, Massachusetts | | |
|---|-----------------------|------------|-----------------------------|---|----------------|---|---|---|--|
| | | | | | | PROJECT LOCATION Walpole, Massach | | | |
| | | | | | | 7 GROUND ELEVATION | | | |
| | | | | | | | | | |
| | | | Hollow Ster | | | | | | |
| LOGGED BY Luke Tulley CHECKED BY | | | | | | | | | |
| NOTE | S | Ι | | | 1 | AFTER DRILLING | | T | |
| о ОЕРТН (#) | SAMPLE TYPE NUMBER | RECOVERY % | BLOW COUNTS (N VALUE) | REMARKS | GRAPHIC LOG | MATERIAL DESCRIPTION | PID (ppm) | WELL DIAGRAM | |
| | S-1 | 50 | 5-7-7-8 (14) | | | Dry, tan/brown, poorly sorted coarse sand and gravel | 0 | Mounted Road Box | |
| | S-2 | 75 | 5-7-7-4 (14) | | | Dry, tan, well sorted medium to fine sand, trace gravel | 0 | - | |
| 10 | S-3 | 80 | 2-3-3-4 (6) | | | Dry, tan, well sorted medium to fine sand, trace gravel | 0 | - | |
| | S-4 | 85 | 3-4-3-4 (7) | | | Dry, tan, well sorted medium to fine sand 16.0 | 0 | - | |
| 20 | S-5 | 25 | 5-8-15-13 (23) | | <u>;</u> ,,,,, | | 0 | - | |
| | S-6 | 12 | _ | | | 23.0 24.0 Dry, tan, poorly sorted medium to fine sand and large angular gravel Boulder | 0 | - | |
| 30 | S-7 | 30 | 9-23 | | \$ | Dry, tan, poorly sorted coarse sand and angular gravel | 0 | | |
| | | | | | | 30.0 | | ■ Bentonite Seal | |
| 40 | S-8 | 50 | 11-16-11- 20 (27) | | | Wet, brown, poorly sorted coarse to fine sand and 41.0 gravel, rocky | 0 | Sand Filter Pack 2" Machine Slotted Well | |
| | S-9 | 15 | - | and Mor - 14 analysis | | Wet, brown, poorly sorted coarse to fine sand and gravel Bottom of hole at 46.0 feet | 0 | Screen 2" PVC Plug | |
| 40 | S-9 | 15 | | Sample taken and submitted for VOC and MCP-14 analysis from 44-46' (RIZ-10-44'-46') | <u>.</u> | Wet, brown, poorly sorted coarse to fine sand and 46.0 gravel Bottom of hole at 46.0 feet. | 0 | 2" PVC F | |

Appendix E

Laboratory Certificates of Analysis



ANALYTICAL REPORT

Lab Number: L0718979

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: Ray Johnson

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 Report Date: 01/03/08

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (200305), NJ (MA935), RI (LAO00065), ME (2006012), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



L0718979

Lab Number:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 **Report Date:** 01/03/08

| Alpha Sample ID | Client ID | Sample Location |
|-----------------|-----------|-----------------|
| L0718979-01 | RIZ-8 | WALPOLE, MA |
| L0718979-02 | RIZ-10 | WALPOLE, MA |
| L0718979-03 | RIZ-9 | WALPOLE, MA |
| L0718979-04 | GHC-6 | WALPOLE, MA |
| L0718979-05 | MW-9 | WALPOLE, MA |
| L0718979-06 | RIZ-3 | WALPOLE, MA |
| L0718979-07 | MW-3 | WALPOLE, MA |

Project Number: 12700058 Report Date: 01/03/08

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| An a | ffirmative response to questions A, B, C & D is required for "Presumptive Certainty" status | |
|------|---|-----|
| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |

| A res | ponse to questions E and F is required for "Presumptive Certainty" status | |
|-------|---|-----|
| Е | Were all QC performance standards and recommendations for the specified method(s) achieved? | NO |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | YES |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L0718979

Lab Number:

Project Name: WALPOLE PARK SOUTH

Case Narrative

The samples were received in accordance with the chain of custody and no significant deviations were encountered during preparation or analysis unless otherwise noted below.

MCP Related Narratives

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

Volatile Organics

In reference to question E:

The WG307181-1/2 LCS/LCSD % recoveries for Dichlorodifluoromethane and the LCS % recovery for 1,4-Dioxane are below the individual acceptance criteria for the compounds, but within the overall method allowances. These are both difficult analytes.

The WG307181-1/2 LCS/LCSD % RPD for 1,4-Dioxane is above the method acceptance criteria.

The WG307363-1/2 LCS/LCSD % recoveries for Dichlorodifluoromethane are below and the LCSD % recoveries for Acetone (a difficult analyte) and 1,4-Dioxane are above the individual acceptance criteria for the compounds, but within the overall method allowances.

Metals

L0718979-01 through -07 were diluted 4x for the analysis of all 6020A analytes due to non-target analyte interference.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative

ALPHA WOODS HOLE LABS

Date: 01/03/08

ORGANICS



VOLATILES



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-01 Date Collected: 12/19/07 11:10

Client ID: RIZ-8

Date Received: 12/21/07 Field Prep: Sample Location: WALPOLE, MA Field Filtered

Matrix: Water Anaytical Method: 60,8260B 12/27/07 17:23 Analytical Date:

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: Date Collected: 12/19/07 11:10

Client ID: RIZ-8 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| p/m-Xylene ND ug/l 1.0 1 o-Xylene ND ug/l 1.0 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Dibromomethane ND ug/l 5.0 1 1,2,3-Trichloropropane ND ug/l 5.0 1 Styrene ND ug/l 5.0 1 Dichlorodifluoromethane ND ug/l 5.0 1 Acetone ND ug/l 5.0 1 Carbon disulfide ND ug/l 5.0 1 2-Butanone ND ug/l 5.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 5.0 1 Tetrahydrofuran ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 2.5 1 1,2-Dibromoethane <th>Parameter</th> <th>Result</th> <th>Qualifier</th> <th>Units</th> <th>RDL</th> <th>Dilution Factor</th> | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------------------------------|--------|-----------|-------|------|-----------------|
| Mothyl tent bulyl ether ND ug/l 1.0 1 p/m-Xylane ND ug/l 1.0 1 c/xylane ND ug/l 1.0 1 c/xylane ND ug/l 1.0 1 c/xylane ND ug/l 5.0 1 Dibromomethane ND ug/l 5.0 1 Syrene ND ug/l 5.0 1 Dichorodifuoromethane ND ug/l 5.0 1 Acetone ND ug/l 5.0 1 Carbon disulfide ND ug/l 5.0 1 2-Burance ND ug/l 5.0 1 2-Burance ND ug/l 5.0 1 4-Mothyl-2-pentanone ND ug/l 5.0 1 2-Burance ND ug/l 5.0 1 1 Carbon disulfide ND ug/l 5.0 1 2-Burance ND ug/l< | Volatile Organics by MCP 8260B | | | | | |
| ND | 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| o-Xylene ND ugfl 1.0 1 cis-1,2-Dichloroethene ND ugfl 0.50 1 Dibromomethane ND ugfl 5.0 1 1,2-3-Trichloropropane ND ugfl 5.0 1 Styrene ND ugfl 5.0 1 Dichlorodiffuoromethane ND ugfl 5.0 1 Acetone ND ugfl 5.0 1 Carbon disulfide ND ugfl 5.0 1 2-Butanone ND ugfl 5.0 1 4-Methyl-2-pentanone ND ugfl 5.0 1 4-Hextonice ND ugfl 5.0 1 2-Busanone ND ugfl 5.0 1 1-Hextonice ND ugfl 5.0 1 1-Hextonice ND ugfl 2.5 1 1-Lacibinomethane ND ugfl 2.5 1 1,2-Dibromoethane | Methyl tert butyl ether | ND | | ug/l | 1.0 | 1 |
| Dibromomethane ND | p/m-Xylene | ND | | ug/l | 1.0 | 1 |
| Dibromomethane ND | o-Xylene | ND | | ug/l | 1.0 | 1 |
| 1,2,3-Trichloropropane ND Ug/l 1,0 1 | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Styrene ND ug/l 1.0 1 Dichlorodifluoromethane ND ug/l 5.0 1 Acetone ND ug/l 5.0 1 Carbon disulfide ND ug/l 5.0 1 2-Butanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 1 Bromobenzene ND ug/l 2.5 1 Bromobenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 Bromobenzene | Dibromomethane | ND | | ug/l | 5.0 | 1 |
| Dichlorodifluoromethane ND ug/l 5.0 1 Acetone ND ug/l 5.0 1 Carbon disulfide ND ug/l 5.0 1 2-Butanone ND ug/l 5.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 2.5 1 1,2-Dichloropropane ND ug/l 2.5 1 1,2-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 1 Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 2.5 1 | 1,2,3-Trichloropropane | ND | | ug/l | 5.0 | 1 |
| Acetone ND ug/l 5.0 1 Carbon disulfide ND ug/l 5.0 1 2-Butanone ND ug/l 5.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Errandoromethane ND ug/l 2.5 1 Bromochloromethane ND ug/l 2.5 1 1-zerbaydrofuran ND ug/l 2.5 1 1-zerbaydrofuran <td>Styrene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>1.0</td> <td>1</td> | Styrene | ND | | ug/l | 1.0 | 1 |
| Carbon disulfide ND ug/l 5.0 1 2-Butanone ND ug/l 5.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 2.5 1 1-2-Dibromorpane ND ug/l 2.5 1 1,2-Dibromorpane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 1 Bromobenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 uer-Butylbenzene ND ug/l 0.50 1 uer-Butylbenzene ND ug/l 0.50 1 uer-Butylbenzene ND ug/l 0.50 1 p-Chl | Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1 |
| 2-Butanone ND ug/l 5.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 2.5 1 1-2-Dibromoethane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.5 1 1,1,1,2-Tetrachloroptrapane ND ug/l 2.5 1 1,1,1,2-Tetrachloroptrapane ND ug/l 2.5 1 Bromobenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 2.5 1 <t< td=""><td>Acetone</td><td>ND</td><td></td><td>ug/l</td><td>5.0</td><td>1</td></t<> | Acetone | ND | | ug/l | 5.0 | 1 |
| 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 10 1 2,2-Dichloropropane ND ug/l 2.5 1 1,2-Dichloropropane ND ug/l 2.0 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 | Carbon disulfide | ND | | ug/l | 5.0 | 1 |
| 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 10 1 2,2-Dichloropropane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.0 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 1 Bromobenzene ND ug/l 2.5 1 Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 -Chlorotoluene ND ug/l 2.5 1 -Chlorotoluene ND ug/l 2.5 1 < | 2-Butanone | ND | | ug/l | 5.0 | 1 |
| Bromochloromethane ND ug/l 2.5 1 | 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1 |
| Tetrahydrofuran ND ug/l 10 1 2,2-Dichloropropane ND ug/l 2.5 1 1,2-Dibromoethane ND ug/l 2.5 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 2.5 1 ec-Butylbenzene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 Hexac | 2-Hexanone | ND | | ug/l | 5.0 | 1 |
| 2,2-Dichloropropane ND ug/l 2,5 1 1,2-Dibromoethane ND ug/l 2,0 1 1,3-Dichloropropane ND ug/l 2,5 1 1,1,1,2-Tetrachloroethane ND ug/l 0,50 1 Bromobenzene ND ug/l 0,50 1 Bromobenzene ND ug/l 0,50 1 sec-Butylbenzene ND ug/l 0,50 1 tert-Butylbenzene ND ug/l 2,5 1 o-Chlorotoluene ND ug/l 2,5 1 p-Chlorotoluene ND ug/l 2,5 1 Hexachlorobutadiene ND ug/l 2,5 1 Hexachlorobutadiene ND ug/l 0,50 1 sopropylbenzene ND ug/l 0,50 1 Naphthalene ND ug/l 0,50 1 n-Propylbenzene ND ug/l 2,5 1 <t< td=""><td>Bromochloromethane</td><td>ND</td><td></td><td>ug/l</td><td>2.5</td><td>1</td></t<> | Bromochloromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromoethane ND ug/l 2.0 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.50 1 sopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 2.5 1 <tr< td=""><td>Tetrahydrofuran</td><td>ND</td><td></td><td>ug/l</td><td>10</td><td>1</td></tr<> | Tetrahydrofuran | ND | | ug/l | 10 | 1 |
| 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 | 2,2-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 2.5 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 <td>1,2-Dibromoethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>2.0</td> <td>1</td> | 1,2-Dibromoethane | ND | | ug/l | 2.0 | 1 |
| ND | 1,3-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| n-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 2.5 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | Bromobenzene | ND | | ug/l | 2.5 | 1 |
| tert-Butylbenzene ND ug/l 2.5 1 o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene ND ug/l 2.5 1 p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 2.5 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene ND ug/l 2.5 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | tert-Butylbenzene | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | o-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | p-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| Sopropylbenzene ND ug/l 0.50 1 | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 1 |
| p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | Hexachlorobutadiene | ND | | ug/l | 0.60 | 1 |
| Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene ND ug/l 2.5 1 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | Naphthalene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trichlorobenzene ND ug/l 2.5 1 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene ND ug/l 2.5 1 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trimethylbenzene ND ug/l 2.5 1 | 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| · | 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| Ethyl ether ND ug/l 2.5 1 | 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| | Ethyl ether | ND | | ug/l | 2.5 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: Date Collected: 12/19/07 11:10

Client ID: RIZ-8 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|-----|------------------------|
| Volatile Organics by MCP 8260B | | | | | |
| | | | | | , |
| Isopropyl Ether | ND | | ug/l | 2.0 | 1 |
| Ethyl-Tert-Butyl-Ether | ND | | ug/l | 2.0 | 1 |
| Tertiary-Amyl Methyl Ether | ND | | ug/l | 2.0 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 1 |

| | Acceptance | | | | | | |
|-----------------------|------------|-----------|----------|--|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | | |
| 1,2-Dichloroethane-d4 | 103 | | 70-130 | | | | |
| Toluene-d8 | 95 | | 70-130 | | | | |
| 4-Bromofluorobenzene | 98 | | 70-130 | | | | |
| Dibromofluoromethane | 105 | | 70-130 | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Lab ID: Date Collected: 12/19/07 12:40

Client ID: RIZ-10 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water
Anaytical Method: 60,8260B
Analytical Date: 12/27/07 18:01

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: Date Collected: 12/19/07 12:40

Client ID: RIZ-10 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| Methyl tert butyl ether | 1.2 | | ug/l | 1.0 | <u>'</u> |
| p/m-Xylene | ND | | ug/l | 1.0 | <u>'</u> |
| o-Xylene | ND | | ug/l | 1.0 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 5.0 | 1 |
| Styrene | ND | | ug/l | 1.0 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 1 |
| Tetrahydrofuran | ND | | ug/l | 10 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.60 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: Date Collected: 12/19/07 12:40

Client ID: RIZ-10 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Result | Qualifier | Units | RDL | Dilution Factor |
|--------|------------------------|----------------|-------------------------|-------------------------------------|
| | | | | |
| ND | | /1 | 2.0 | 1 |
| | | • | | 1 |
| | | | | 1 |
| | | <u> </u> | | 1 |
| | Result ND ND ND ND | ND ND ND | ND ug/l ND ug/l ND ug/l | ND ug/l 2.0 ND ug/l 2.0 ND ug/l 2.0 |

| | Acceptance | | | | | | |
|-----------------------|------------|-----------|----------|--|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | | |
| 1,2-Dichloroethane-d4 | 97 | | 70-130 | | | | |
| Toluene-d8 | 96 | | 70-130 | | | | |
| 4-Bromofluorobenzene | 99 | | 70-130 | | | | |
| Dibromofluoromethane | 99 | | 70-130 | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-03 Date Collected: 12/19/07 13:40

Client ID: RIZ-9

Date Received: 12/21/07 Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water Anaytical Method: 60,8260B 12/27/07 18:40 Analytical Date:

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-03 Date Collected: 12/19/07 13:40

Client ID: RIZ-9 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 1.0 | 1 |
| p/m-Xylene | ND | | ug/l | 1.0 | 1 |
| o-Xylene | ND | | ug/l | 1.0 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 5.0 | 1 |
| Styrene | ND | | ug/l | 1.0 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 1 |
| Tetrahydrofuran | ND | | ug/l | 10 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.60 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-03 Date Collected: 12/19/07 13:40

Client ID: RIZ-9 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|-----|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Isopropyl Ether | ND | | ug/l | 2.0 | 1 |
| Ethyl-Tert-Butyl-Ether | ND | | ug/l | 2.0 | 1 |
| Tertiary-Amyl Methyl Ether | ND | | ug/l | 2.0 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|-----------------------|------------|-----------|------------------------|--|
| 1,2-Dichloroethane-d4 | 100 | | 70-130 | |
| Toluene-d8 | 96 | | 70-130 | |
| 4-Bromofluorobenzene | 99 | | 70-130 | |
| Dibromofluoromethane | 102 | | 70-130 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-04 Date Collected: 12/19/07 16:00

Client ID: GHC-6 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water
Anaytical Method: 60,8260B
Analytical Date: 12/27/07 19:18

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-04 Date Collected: 12/19/07 16:00

Client ID: GHC-6 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

Parameter Result Qualifier Units **RDL Dilution Factor** Volatile Organics by MCP 8260B 1,4-Dichlorobenzene ND 2.5 ug/l 1 ND 1.0 1 Methyl tert butyl ether ug/l ND p/m-Xylene ug/l 1.0 1 o-Xylene ND ug/l 1.0 1 ND cis-1,2-Dichloroethene ug/l 0.50 1 Dibromomethane ND 5.0 ug/l 1 1,2,3-Trichloropropane ND ug/l 5.0 1 ND 1.0 1 Styrene ug/l Dichlorodifluoromethane ND 5.0 1 ug/l ND 5.0 Acetone ug/l 1 Carbon disulfide ND 5.0 ug/l 1 2-Butanone ND 5.0 1 ug/l 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 10 1 ND 1 2,2-Dichloropropane ug/l 2.5 1,2-Dibromoethane ND ug/l 2.0 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 ND 0.50 sec-Butylbenzene ug/l 1 ND 2.5 1 tert-Butylbenzene ug/l ND o-Chlorotoluene ug/l 2.5 1 p-Chlorotoluene ND 2.5 1 ug/l ND 1,2-Dibromo-3-chloropropane ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND 0.50 1 ug/l Naphthalene ND ug/l 2.5 1 ND n-Propylbenzene ug/l 0.50 1 1,2,3-Trichlorobenzene ND 2.5 ug/l 1 1,2,4-Trichlorobenzene ND 2.5 1 ug/l 1,3,5-Trimethylbenzene ND ug/l 2.5 1 ND 1,2,4-Trimethylbenzene 2.5 1 ug/l Ethyl ether ND 2.5 1 ug/l



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-04 Date Collected: 12/19/07 16:00

Client ID: GHC-6 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|-----|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Isopropyl Ether | ND | | ug/l | 2.0 | 1 |
| Ethyl-Tert-Butyl-Ether | ND | | ug/l | 2.0 | 1 |
| Tertiary-Amyl Methyl Ether | ND | | ug/l | 2.0 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 1 |

| | Acceptance | | | | | | |
|-----------------------|------------|-----------|----------|--|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | | |
| 1,2-Dichloroethane-d4 | 96 | | 70-130 | | | | |
| Toluene-d8 | 97 | | 70-130 | | | | |
| 4-Bromofluorobenzene | 99 | | 70-130 | | | | |
| Dibromofluoromethane | 102 | | 70-130 | | | | |



12/20/07 09:00

Field Filtered

12/21/07

Date Collected:

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-05

Client ID: MW-9

Sample Location: WALPOLE, MA

Matrix: Water
Anaytical Method: 60,8260B
Analytical Date: 12/27/07 19:57

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-05 Date Collected: 12/20/07 09:00

Client ID: MW-9 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| Methyl tert butyl ether | 5.1 | | ug/l | 1.0 | 1 |
| p/m-Xylene | ND | | ug/l | 1.0 | 1 |
| o-Xylene | ND | | ug/l | 1.0 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 5.0 | 1 |
| Styrene | ND | | ug/l | 1.0 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 1 |
| Tetrahydrofuran | ND | | ug/l | 10 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.60 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 1 |



1

250

ug/l

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-05 Date Collected: 12/20/07 09:00

Client ID: MW-9 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

Parameter Qualifier Units **RDL Dilution Factor** Result Volatile Organics by MCP 8260B Isopropyl Ether ND 2.0 ug/l 1 Ethyl-Tert-Butyl-Ether ND ug/l 2.0 1 Tertiary-Amyl Methyl Ether ND ug/l 2.0 1

| | Acceptance | | | | | |
|-----------------------|------------|-----------|-----------------|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | |
| 1,2-Dichloroethane-d4 | 98 | | 70-130 | | | |
| Toluene-d8 | 95 | | 70-130 | | | |
| 4-Bromofluorobenzene | 98 | | 70-130 | | | |
| Dibromofluoromethane | 101 | | 70-130 | | | |

ND



1,4-Dioxane

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-06 Date Collected: 12/20/07 10:30

Client ID: RIZ-3

Date Received: 12/21/07 Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water Anaytical Method: 60,8260B Analytical Date: 12/27/07 20:36

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-06 Date Collected: 12/20/07 10:30

Client ID: RIZ-3 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 1.0 | 1 |
| p/m-Xylene | ND | | ug/l | 1.0 | 1 |
| o-Xylene | ND | | ug/l | 1.0 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 5.0 | 1 |
| Styrene | ND | | ug/l | 1.0 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 1 |
| Tetrahydrofuran | ND | | ug/l | 10 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.60 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-06 Date Collected: 12/20/07 10:30

Client ID: RIZ-3 Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|-----|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Isopropyl Ether | ND | | ug/l | 2.0 | 1 |
| Ethyl-Tert-Butyl-Ether | ND | | ug/l | 2.0 | 1 |
| Tertiary-Amyl Methyl Ether | ND | | ug/l | 2.0 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|-----------------------|------------|-----------|------------------------|--|
| 1,2-Dichloroethane-d4 | 101 | | 70-130 | |
| Toluene-d8 | 95 | | 70-130 | |
| 4-Bromofluorobenzene | 97 | | 70-130 | |
| Dibromofluoromethane | 100 | | 70-130 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-07 Date Collected: 12/20/07 11:25

Client ID: MW-3

Date Received: 12/21/07 Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water Anaytical Method: 60,8260B Analytical Date: 12/28/07 21:58

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Methylene chloride | ND | | ug/l | 5.0 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.75 | 1 |
| Chloroform | ND | | ug/l | 0.75 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.8 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.75 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.75 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 1 |
| Bromomethane | ND | | ug/l | 1.0 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 1 |
| Chloroethane | ND | | ug/l | 1.0 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.75 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-07 Date Collected: 12/20/07 11:25

Client ID: MW-3 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

Parameter Result Qualifier Units **RDL Dilution Factor** Volatile Organics by MCP 8260B 1,4-Dichlorobenzene ND 2.5 ug/l 1 1.7 1.0 1 Methyl tert butyl ether ug/l ND p/m-Xylene ug/l 1.0 1 o-Xylene ND ug/l 1.0 1 ND cis-1,2-Dichloroethene ug/l 0.50 1 Dibromomethane ND 5.0 ug/l 1 1,2,3-Trichloropropane ND ug/l 5.0 1 ND 1.0 1 Styrene ug/l Dichlorodifluoromethane ND 5.0 1 ug/l ND 5.0 Acetone ug/l 1 Carbon disulfide ND 5.0 ug/l 1 2-Butanone ND 5.0 1 ug/l 4-Methyl-2-pentanone ND ug/l 5.0 1 2-Hexanone ND ug/l 5.0 1 Bromochloromethane ND ug/l 2.5 1 Tetrahydrofuran ND ug/l 10 1 ND 1 2,2-Dichloropropane ug/l 2.5 1,2-Dibromoethane ND ug/l 2.0 1 1,3-Dichloropropane ND ug/l 2.5 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l Bromobenzene ND ug/l 2.5 1 n-Butylbenzene ND ug/l 0.50 1 ND 0.50 sec-Butylbenzene ug/l 1 ND 2.5 1 tert-Butylbenzene ug/l ND o-Chlorotoluene ug/l 2.5 1 p-Chlorotoluene ND 2.5 1 ug/l ND 1,2-Dibromo-3-chloropropane ug/l 2.5 1 Hexachlorobutadiene ND ug/l 0.60 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND 0.50 1 ug/l Naphthalene ND ug/l 2.5 1 n-Propylbenzene ND ug/l 0.50 1 ND 2.5 1,2,3-Trichlorobenzene ug/l 1 1,2,4-Trichlorobenzene ND 2.5 1 ug/l 1,3,5-Trimethylbenzene ND ug/l 2.5 1 ND 1,2,4-Trimethylbenzene 2.5 1 ug/l Ethyl ether ND 2.5 1 ug/l



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-07 Date Collected: 12/20/07 11:25

Client ID: MW-3 Date Received: 12/21/07
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------|--------|-----------|-------|-----|-----------------|
| Volatile Organics by MCP 8260B | | | | | |
| Isopropyl Ether | ND | | ug/l | 2.0 | 1 |
| Ethyl-Tert-Butyl-Ether | ND | | ug/l | 2.0 | 1 |
| Tertiary-Amyl Methyl Ether | ND | | ug/l | 2.0 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 1 |

| | Acceptance | | | | | | | |
|-----------------------|------------|-----------|-----------------|--|--|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | | | |
| 1,2-Dichloroethane-d4 | 107 | | 70-130 | | | | | |
| Toluene-d8 | 94 | | 70-130 | | | | | |
| 4-Bromofluorobenzene | 97 | | 70-130 | | | | | |
| Dibromofluoromethane | 107 | | 70-130 | | | | | |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/27/07 12:14

| Result | Qualifie | r U | nits | RDL |
|------------------|--|--|---|--|
| B for sample(s): | 01-06 | Batch: | WG: | 307181-3 |
| ND | | ı | ug/l | 5.0 |
| ND | | | ug/l | 0.75 |
| ND | | | ug/l | 0.75 |
| ND | | | ug/l | 0.50 |
| ND | | | ug/l | 1.8 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.75 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 2.5 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 2.5 |
| ND | | ı | ug/l | 2.0 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.75 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 2.5 |
| ND | | ı | ug/l | 1.0 |
| ND | | ı | ug/l | 1.0 |
| ND | | ı | ug/l | 1.0 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 0.75 |
| ND | | ı | ug/l | 0.50 |
| ND | | ı | ug/l | 2.5 |
| ND | | ı | ug/l | 2.5 |
| ND | | ı | ug/l | 2.5 |
| | ND ND ND ND ND ND ND ND ND ND ND ND ND N | ND ND ND ND ND ND ND ND ND ND ND ND ND N | DB for sample(s): 01-06 Batch: ND ND ND ND ND ND ND ND ND N | ND ug/l ND |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/27/07 12:14

| arameter | Result | Qualifie | r U | nits | RDL |
|-----------------------------|----------------------|----------|--------|------|----------|
| olatile Organics by MCP | 8260B for sample(s): | 01-06 | Batch: | WG: | 307181-3 |
| Methyl tert butyl ether | ND | | ı | ug/l | 1.0 |
| p/m-Xylene | ND | | | ug/l | 1.0 |
| o-Xylene | ND | | | ug/l | 1.0 |
| cis-1,2-Dichloroethene | ND | | | ug/l | 0.50 |
| Dibromomethane | ND | | | ug/l | 5.0 |
| 1,2,3-Trichloropropane | ND | | (| ug/l | 5.0 |
| Styrene | ND | | ı | ug/l | 1.0 |
| Dichlorodifluoromethane | ND | | ı | ug/l | 5.0 |
| Acetone | ND | | ı | ug/l | 5.0 |
| Carbon disulfide | ND | | ı | ug/l | 5.0 |
| 2-Butanone | ND | | ı | ug/l | 5.0 |
| 4-Methyl-2-pentanone | ND | | ı | ug/l | 5.0 |
| 2-Hexanone | ND | | ı | ug/l | 5.0 |
| Bromochloromethane | ND | | ı | ug/l | 2.5 |
| Tetrahydrofuran | ND | | ı | ug/l | 10 |
| 2,2-Dichloropropane | ND | | (| ug/l | 2.5 |
| 1,2-Dibromoethane | ND | | ı | ug/l | 2.0 |
| 1,3-Dichloropropane | ND | | ı | ug/l | 2.5 |
| 1,1,1,2-Tetrachloroethane | ND | | ı | ug/l | 0.50 |
| Bromobenzene | ND | | ı | ug/l | 2.5 |
| n-Butylbenzene | ND | | ı | ug/l | 0.50 |
| sec-Butylbenzene | ND | | ı | ug/l | 0.50 |
| tert-Butylbenzene | ND | | ı | ug/l | 2.5 |
| o-Chlorotoluene | ND | | ı | ug/l | 2.5 |
| p-Chlorotoluene | ND | | ı | ug/l | 2.5 |
| 1,2-Dibromo-3-chloropropane | ND | | ı | ug/l | 2.5 |
| Hexachlorobutadiene | ND | | | ug/l | 0.60 |
| Isopropylbenzene | ND | | | ug/l | 0.50 |
| p-Isopropyltoluene | ND | | ı | ug/l | 0.50 |
| Naphthalene | ND | | | ug/l | 2.5 |
| n-Propylbenzene | ND | | ı | ug/l | 0.50 |
| | | | | | |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/27/07 12:14

| olatile Organics by MCP 8260B for sample(s): 01-06 Batch: WG307181-3 |
|--|
| 1,2,3-Trichlorobenzene ND ug/l 2.5 |
| 1,2,4-Trichlorobenzene ND ug/l 2.5 |
| 1,3,5-Trimethylbenzene ND ug/l 2.5 |
| 1,2,4-Trimethylbenzene ND ug/l 2.5 |
| Ethyl ether ND ug/l 2.5 |
| Isopropyl Ether ND ug/l 2.0 |
| Ethyl-Tert-Butyl-Ether ND ug/l 2.0 |
| Tertiary-Amyl Methyl Ether ND ug/l 2.0 |
| 1,4-Dioxane ND ug/l 250 |

| | | Acceptance | | | | | | |
|-----------------------|-----------|------------|----------|--|--|--|--|--|
| Surrogate | %Recovery | Qualifier | Criteria | | | | | |
| 1,2-Dichloroethane-d4 | 104 | | 70-130 | | | | | |
| Toluene-d8 | 94 | | 70-130 | | | | | |
| 4-Bromofluorobenzene | 98 | | 70-130 | | | | | |
| Dibromofluoromethane | 107 | | 70-130 | | | | | |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/28/07 13:00

| arameter | Result | Qual | lifier | Units | RDL |
|-----------------------------|-------------------|------|--------|-------|-------|
| olatile Organics by MCP 826 | 0B for sample(s): | 07 | Batch: | WG307 | 363-3 |
| Methylene chloride | ND | | | ug/l | 5.0 |
| 1,1-Dichloroethane | ND | | | ug/l | 0.75 |
| Chloroform | ND | | | ug/l | 0.75 |
| Carbon tetrachloride | ND | | | ug/l | 0.50 |
| 1,2-Dichloropropane | ND | | | ug/l | 1.8 |
| Dibromochloromethane | ND | | | ug/l | 0.50 |
| 1,1,2-Trichloroethane | ND | | | ug/l | 0.75 |
| Tetrachloroethene | ND | | | ug/l | 0.50 |
| Chlorobenzene | ND | | | ug/l | 0.50 |
| Trichlorofluoromethane | ND | | | ug/l | 2.5 |
| 1,2-Dichloroethane | ND | | | ug/l | 0.50 |
| 1,1,1-Trichloroethane | ND | | | ug/l | 0.50 |
| Bromodichloromethane | ND | | | ug/l | 0.50 |
| trans-1,3-Dichloropropene | ND | | | ug/l | 0.50 |
| cis-1,3-Dichloropropene | ND | | | ug/l | 0.50 |
| 1,1-Dichloropropene | ND | | | ug/l | 2.5 |
| Bromoform | ND | | | ug/l | 2.0 |
| 1,1,2,2-Tetrachloroethane | ND | | | ug/l | 0.50 |
| Benzene | ND | | | ug/l | 0.50 |
| Toluene | ND | | | ug/l | 0.75 |
| Ethylbenzene | ND | | | ug/l | 0.50 |
| Chloromethane | ND | | | ug/l | 2.5 |
| Bromomethane | ND | | | ug/l | 1.0 |
| Vinyl chloride | ND | | | ug/l | 1.0 |
| Chloroethane | ND | | | ug/l | 1.0 |
| 1,1-Dichloroethene | ND | | | ug/l | 0.50 |
| trans-1,2-Dichloroethene | ND | | | ug/l | 0.75 |
| Trichloroethene | ND | | | ug/l | 0.50 |
| 1,2-Dichlorobenzene | ND | | | ug/l | 2.5 |
| 1,3-Dichlorobenzene | ND | | | ug/l | 2.5 |
| 1,4-Dichlorobenzene | ND | | | ug/l | 2.5 |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/28/07 13:00

| Parameter | Result | Qua | lifier | Units | RDL |
|----------------------------------|---------------|-----|--------|--------|-------|
| olatile Organics by MCP 8260B fc | or sample(s): | 07 | Batch: | WG3073 | 363-3 |
| Methyl tert butyl ether | ND | | | ug/l | 1.0 |
| p/m-Xylene | ND | | | ug/l | 1.0 |
| o-Xylene | ND | | | ug/l | 1.0 |
| cis-1,2-Dichloroethene | ND | | | ug/l | 0.50 |
| Dibromomethane | ND | | | ug/l | 5.0 |
| 1,2,3-Trichloropropane | ND | | | ug/l | 5.0 |
| Styrene | ND | | | ug/l | 1.0 |
| Dichlorodifluoromethane | ND | | | ug/l | 5.0 |
| Acetone | ND | | | ug/l | 5.0 |
| Carbon disulfide | ND | | | ug/l | 5.0 |
| 2-Butanone | ND | | | ug/l | 5.0 |
| 4-Methyl-2-pentanone | ND | | | ug/l | 5.0 |
| 2-Hexanone | ND | | | ug/l | 5.0 |
| Bromochloromethane | ND | | | ug/l | 2.5 |
| Tetrahydrofuran | ND | | | ug/l | 10 |
| 2,2-Dichloropropane | ND | | | ug/l | 2.5 |
| 1,2-Dibromoethane | ND | | | ug/l | 2.0 |
| 1,3-Dichloropropane | ND | | | ug/l | 2.5 |
| 1,1,1,2-Tetrachloroethane | ND | | | ug/l | 0.50 |
| Bromobenzene | ND | | | ug/l | 2.5 |
| n-Butylbenzene | ND | | | ug/l | 0.50 |
| sec-Butylbenzene | ND | | | ug/l | 0.50 |
| tert-Butylbenzene | ND | | | ug/l | 2.5 |
| o-Chlorotoluene | ND | | | ug/l | 2.5 |
| p-Chlorotoluene | ND | | | ug/l | 2.5 |
| 1,2-Dibromo-3-chloropropane | ND | | | ug/l | 2.5 |
| Hexachlorobutadiene | ND | | | ug/l | 0.60 |
| Isopropylbenzene | ND | | | ug/l | 0.50 |
| p-Isopropyltoluene | ND | | | ug/l | 0.50 |
| Naphthalene | ND | | | ug/l | 2.5 |
| n-Propylbenzene | ND | | | ug/l | 0.50 |
| | | | | | |



Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

Analytical Method: 60,8260B Analytical Date: 12/28/07 13:00

| arameter | Result | Qual | ifier | Units | RDL |
|-----------------------------------|------------|------|--------|--------|-------|
| olatile Organics by MCP 8260B for | sample(s): | 07 | Batch: | WG3073 | 363-3 |
| 1,2,3-Trichlorobenzene | ND | | | ug/l | 2.5 |
| 1,2,4-Trichlorobenzene | ND | | | ug/l | 2.5 |
| 1,3,5-Trimethylbenzene | ND | | | ug/l | 2.5 |
| 1,2,4-Trimethylbenzene | ND | | | ug/l | 2.5 |
| Ethyl ether | ND | | | ug/l | 2.5 |
| Isopropyl Ether | ND | | | ug/l | 2.0 |
| Ethyl-Tert-Butyl-Ether | ND | | | ug/l | 2.0 |
| Tertiary-Amyl Methyl Ether | ND | | | ug/l | 2.0 |
| 1,4-Dioxane | ND | | | ug/l | 250 |

| | Acceptance | | | | | | |
|-----------------------|------------|--------------------|---|--|--|--|--|
| Surrogate | %Recovery | Qualifier Criteria | _ | | | | |
| 1,2-Dichloroethane-d4 | 108 | 70-130 | | | | | |
| Toluene-d8 | 95 | 70-130 | | | | | |
| 4-Bromofluorobenzene | 94 | 70-130 | | | | | |
| Dibromofluoromethane | 107 | 70-130 | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

| Parameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---|------------------|-------------------|---------------------|-----|------------|
| /olatile Organics by MCP 8260B Associated | sample(s): 01-06 | Batch: WG307181 | -1 WG307181-2 | | |
| Methylene chloride | 99 | 101 | 70-130 | 2 | 25 |
| 1,1-Dichloroethane | 100 | 100 | 70-130 | 0 | 25 |
| Chloroform | 102 | 104 | 70-130 | 2 | 25 |
| Carbon tetrachloride | 89 | 98 | 70-130 | 10 | 25 |
| 1,2-Dichloropropane | 95 | 97 | 70-130 | 2 | 25 |
| Dibromochloromethane | 86 | 92 | 70-130 | 7 | 25 |
| 1,1,2-Trichloroethane | 91 | 87 | 70-130 | 4 | 25 |
| Tetrachloroethene | 99 | 103 | 70-130 | 4 | 25 |
| Chlorobenzene | 96 | 98 | 70-130 | 2 | 25 |
| Trichlorofluoromethane | 105 | 108 | 70-130 | 3 | 25 |
| 1,2-Dichloroethane | 97 | 100 | 70-130 | 3 | 25 |
| 1,1,1-Trichloroethane | 99 | 102 | 70-130 | 3 | 25 |
| Bromodichloromethane | 93 | 98 | 70-130 | 5 | 25 |
| trans-1,3-Dichloropropene | 84 | 86 | 70-130 | 2 | 25 |
| cis-1,3-Dichloropropene | 94 | 95 | 70-130 | 1 | 25 |
| 1,1-Dichloropropene | 99 | 102 | 70-130 | 3 | 25 |
| Bromoform | 82 | 89 | 70-130 | 8 | 50 |
| 1,1,2,2-Tetrachloroethane | 84 | 85 | 70-130 | 1 | 25 |
| Benzene | 97 | 99 | 70-130 | 2 | 25 |
| Toluene | 95 | 97 | 70-130 | 2 | 25 |
| Ethylbenzene | 97 | 98 | 70-130 | 1 | 25 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

| Parameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|-------------------------------------|--------------------------|-------------------|---------------------|-----|------------|
| Volatile Organics by MCP 8260B Asso | ociated sample(s): 01-06 | Batch: WG307181-1 | WG307181-2 | | |
| Chloromethane | 78 | 83 | 70-130 | 6 | 50 |
| Bromomethane | 91 | 89 | 70-130 | 2 | 50 |
| Vinyl chloride | 83 | 88 | 70-130 | 6 | 25 |
| Chloroethane | 111 | 106 | 70-130 | 5 | 25 |
| 1,1-Dichloroethene | 102 | 104 | 70-130 | 2 | 25 |
| trans-1,2-Dichloroethene | 96 | 99 | 70-130 | 3 | 25 |
| Trichloroethene | 96 | 100 | 70-130 | 4 | 25 |
| 1,2-Dichlorobenzene | 99 | 97 | 70-130 | 2 | 25 |
| 1,3-Dichlorobenzene | 98 | 101 | 70-130 | 3 | 25 |
| 1,4-Dichlorobenzene | 98 | 98 | 70-130 | 0 | 25 |
| Methyl tert butyl ether | 100 | 98 | 70-130 | 2 | 25 |
| p/m-Xylene | 101 | 104 | 70-130 | 3 | 25 |
| o-Xylene | 101 | 102 | 70-130 | 1 | 25 |
| cis-1,2-Dichloroethene | 97 | 100 | 70-130 | 3 | 25 |
| Dibromomethane | 93 | 94 | 70-130 | 1 | 25 |
| 1,2,3-Trichloropropane | 86 | 88 | 70-130 | 2 | 25 |
| Styrene | 100 | 102 | 70-130 | 2 | 25 |
| Dichlorodifluoromethane | 60 | 65 | 70-130 | 8 | 50 |
| Acetone | 95 | 92 | 70-130 | 3 | 50 |
| Carbon disulfide | 90 | 91 | 70-130 | 1 | 25 |
| 2-Butanone | 87 | 90 | 70-130 | 3 | 50 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|--|--------------------|-------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B Associated | d sample(s): 01-06 | Batch: WG307181- | 1 WG307181-2 | | |
| 4-Methyl-2-pentanone | 93 | 99 | 70-130 | 6 | 50 |
| 2-Hexanone | 84 | 86 | 70-130 | 2 | 50 |
| Bromochloromethane | 100 | 106 | 70-130 | 6 | 25 |
| Tetrahydrofuran | 91 | 94 | 70-130 | 3 | 25 |
| 2,2-Dichloropropane | 100 | 105 | 70-130 | 5 | 50 |
| 1,2-Dibromoethane | 90 | 91 | 70-130 | 1 | 25 |
| 1,3-Dichloropropane | 88 | 90 | 70-130 | 2 | 25 |
| 1,1,1,2-Tetrachloroethane | 88 | 93 | 70-130 | 6 | 25 |
| Bromobenzene | 98 | 97 | 70-130 | 1 | 25 |
| n-Butylbenzene | 95 | 96 | 70-130 | 1 | 25 |
| sec-Butylbenzene | 96 | 98 | 70-130 | 2 | 25 |
| tert-Butylbenzene | 98 | 98 | 70-130 | 0 | 25 |
| o-Chlorotoluene | 94 | 94 | 70-130 | 0 | 25 |
| p-Chlorotoluene | 95 | 96 | 70-130 | 1 | 25 |
| 1,2-Dibromo-3-chloropropane | 74 | 87 | 70-130 | 16 | 50 |
| Hexachlorobutadiene | 96 | 101 | 70-130 | 5 | 25 |
| Isopropylbenzene | 103 | 106 | 70-130 | 3 | 25 |
| p-Isopropyltoluene | 102 | 102 | 70-130 | 0 | 25 |
| Naphthalene | 90 | 97 | 70-130 | 7 | 25 |
| n-Propylbenzene | 96 | 96 | 70-130 | 0 | 25 |
| 1,2,3-Trichlorobenzene | 100 | 105 | 70-130 | 5 | 25 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L0718979

01/03/08

Report Date:

| arameter | LCS %Recover | у | % | LCSD Recovery | %Recovery Limits | RPD | RPD Limits |
|---------------------------------|-----------------------|-------|--------|------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B A | Associated sample(s): | 01-06 | Batch: | WG307181-1 | WG307181-2 | | |
| 1,2,4-Trichlorobenzene | 103 | | | 106 | 70-130 | 3 | 25 |
| 1,3,5-Trimethylbenzene | 97 | | | 96 | 70-130 | 1 | 25 |
| 1,2,4-Trimethylbenzene | 97 | | | 97 | 70-130 | 0 | 25 |
| Ethyl ether | 103 | | | 103 | 70-130 | 0 | 25 |
| Isopropyl Ether | 100 | | | 100 | 70-130 | 0 | 25 |
| Ethyl-Tert-Butyl-Ether | 101 | | | 94 | 70-130 | 7 | 25 |
| Tertiary-Amyl Methyl Ether | 96 | | | 100 | 70-130 | 4 | 25 |
| 1,4-Dioxane | 53 | | | 113 | 70-130 | 72 | 50 |

| Surrogate | LCS %Recovery Qualifier | LCSD %Recovery Qualifier | Acceptance Criteria |
|-----------------------|----------------------------|-----------------------------|------------------------|
| 1,2-Dichloroethane-d4 | 99 | 100 | 70-130 |
| Toluene-d8 | 98 | 95 | 70-130 |
| 4-Bromofluorobenzene | 97 | 96 | 70-130 |
| Dibromofluoromethane | 102 | 102 | 70-130 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|--|--------------------|-------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B Associated | d sample(s): 07 Ba | tch: WG307363-1 | WG307363-2 | | |
| Methylene chloride | 99 | 100 | 70-130 | 1 | 25 |
| 1,1-Dichloroethane | 103 | 99 | 70-130 | 4 | 25 |
| Chloroform | 108 | 105 | 70-130 | 3 | 25 |
| Carbon tetrachloride | 111 | 101 | 70-130 | 9 | 25 |
| 1,2-Dichloropropane | 94 | 91 | 70-130 | 3 | 25 |
| Dibromochloromethane | 94 | 92 | 70-130 | 2 | 25 |
| 1,1,2-Trichloroethane | 88 | 92 | 70-130 | 4 | 25 |
| Tetrachloroethene | 104 | 98 | 70-130 | 6 | 25 |
| Chlorobenzene | 96 | 91 | 70-130 | 5 | 25 |
| Trichlorofluoromethane | 118 | 110 | 70-130 | 7 | 25 |
| 1,2-Dichloroethane | 104 | 109 | 70-130 | 5 | 25 |
| 1,1,1-Trichloroethane | 111 | 107 | 70-130 | 4 | 25 |
| Bromodichloromethane | 102 | 100 | 70-130 | 2 | 25 |
| trans-1,3-Dichloropropene | 88 | 86 | 70-130 | 2 | 25 |
| cis-1,3-Dichloropropene | 100 | 98 | 70-130 | 2 | 25 |
| 1,1-Dichloropropene | 108 | 101 | 70-130 | 7 | 25 |
| Bromoform | 95 | 93 | 70-130 | 2 | 50 |
| 1,1,2,2-Tetrachloroethane | 82 | 84 | 70-130 | 2 | 25 |
| Benzene | 100 | 95 | 70-130 | 5 | 25 |
| Toluene | 94 | 88 | 70-130 | 7 | 25 |
| Ethylbenzene | 97 | 92 | 70-130 | 5 | 25 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

Report Date: 01/03/08

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|--|-------------------|-------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B Associated | sample(s): 07 Bar | tch: WG307363-1 | WG307363-2 | | |
| Chloromethane | 80 | 76 | 70-130 | 5 | 50 |
| Bromomethane | 82 | 75 | 70-130 | 9 | 50 |
| Vinyl chloride | 88 | 83 | 70-130 | 6 | 25 |
| Chloroethane | 105 | 100 | 70-130 | 5 | 25 |
| 1,1-Dichloroethene | 111 | 104 | 70-130 | 7 | 25 |
| trans-1,2-Dichloroethene | 102 | 97 | 70-130 | 5 | 25 |
| Trichloroethene | 103 | 97 | 70-130 | 6 | 25 |
| 1,2-Dichlorobenzene | 94 | 93 | 70-130 | 1 | 25 |
| 1,3-Dichlorobenzene | 96 | 91 | 70-130 | 5 | 25 |
| 1,4-Dichlorobenzene | 97 | 93 | 70-130 | 4 | 25 |
| Methyl tert butyl ether | 106 | 117 | 70-130 | 10 | 25 |
| p/m-Xylene | 102 | 96 | 70-130 | 6 | 25 |
| o-Xylene | 101 | 94 | 70-130 | 7 | 25 |
| cis-1,2-Dichloroethene | 102 | 106 | 70-130 | 4 | 25 |
| Dibromomethane | 100 | 105 | 70-130 | 5 | 25 |
| 1,2,3-Trichloropropane | 82 | 92 | 70-130 | 11 | 25 |
| Styrene | 99 | 95 | 70-130 | 4 | 25 |
| Dichlorodifluoromethane | 66 | 63 | 70-130 | 5 | 50 |
| Acetone | 118 | 147 | 70-130 | 22 | 50 |
| Carbon disulfide | 89 | 87 | 70-130 | 2 | 25 |
| 2-Butanone | 100 | 112 | 70-130 | 11 | 50 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

Report Date: 01/03/08

| arameter | LCS %Recovery | | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---|------------------|--------|-------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B Associate | d sample(s): 07 | Batch: | WG307363-1 | WG307363-2 | | |
| 4-Methyl-2-pentanone | 101 | | 116 | 70-130 | 14 | 50 |
| 2-Hexanone | 85 | | 98 | 70-130 | 14 | 50 |
| Bromochloromethane | 103 | | 108 | 70-130 | 5 | 25 |
| Tetrahydrofuran | 99 | | 114 | 70-130 | 14 | 25 |
| 2,2-Dichloropropane | 111 | | 108 | 70-130 | 3 | 50 |
| 1,2-Dibromoethane | 92 | | 93 | 70-130 | 1 | 25 |
| 1,3-Dichloropropane | 88 | | 93 | 70-130 | 6 | 25 |
| 1,1,1,2-Tetrachloroethane | 95 | | 91 | 70-130 | 4 | 25 |
| Bromobenzene | 97 | | 92 | 70-130 | 5 | 25 |
| n-Butylbenzene | 98 | | 91 | 70-130 | 7 | 25 |
| sec-Butylbenzene | 98 | | 91 | 70-130 | 7 | 25 |
| tert-Butylbenzene | 100 | | 93 | 70-130 | 7 | 25 |
| o-Chlorotoluene | 93 | | 87 | 70-130 | 7 | 25 |
| p-Chlorotoluene | 94 | | 89 | 70-130 | 5 | 25 |
| 1,2-Dibromo-3-chloropropane | 83 | | 85 | 70-130 | 2 | 50 |
| Hexachlorobutadiene | 106 | | 101 | 70-130 | 5 | 25 |
| Isopropylbenzene | 106 | | 99 | 70-130 | 7 | 25 |
| p-Isopropyltoluene | 104 | | 97 | 70-130 | 7 | 25 |
| Naphthalene | 93 | | 96 | 70-130 | 3 | 25 |
| n-Propylbenzene | 96 | | 89 | 70-130 | 8 | 25 |
| 1,2,3-Trichlorobenzene | 100 | | 104 | 70-130 | 4 | 25 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L0718979

Report Date:

| e: 01/03/08 |
|-------------|
| |

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---|-------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by MCP 8260B Associate | ed sample(s): 07 Batch: | WG307363-1 | WG307363-2 | | |
| 1,2,4-Trichlorobenzene | 102 | 100 | 70-130 | 2 | 25 |
| 1,3,5-Trimethylbenzene | 98 | 90 | 70-130 | 9 | 25 |
| 1,2,4-Trimethylbenzene | 95 | 90 | 70-130 | 5 | 25 |
| Ethyl ether | 103 | 116 | 70-130 | 12 | 25 |
| Isopropyl Ether | 101 | 104 | 70-130 | 3 | 25 |
| Ethyl-Tert-Butyl-Ether | 104 | 107 | 70-130 | 3 | 25 |
| Tertiary-Amyl Methyl Ether | 105 | 114 | 70-130 | 8 | 25 |
| 1,4-Dioxane | 130 | 144 | 70-130 | 10 | 50 |

| Surrogate | LCS %Recovery Qualifier | LCSD %Recovery Qualifier | Acceptance Criteria |
|-----------------------|----------------------------|-----------------------------|------------------------|
| 1,2-Dichloroethane-d4 | 106 | 109 | 70-130 |
| Toluene-d8 | 93 | 95 | 70-130 |
| 4-Bromofluorobenzene | 97 | 96 | 70-130 |
| Dibromofluoromethane | 109 | 109 | 70-130 |



METALS



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-01

Client ID: RIZ-8

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 12/19/07 11:10
Date Received: 12/21/07
Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|--------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | MCP 60 | 00/7000 seri | es | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0508 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:31 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | 0.0048 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | вм |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | ND | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 22:50 | EPA 3005A | 64,6020A | ВМ |



12/19/07 12:40

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Date Collected:

Lab ID: L0718979-02

Client ID: Date Received: 12/21/07

Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|--------|--------|-----------------|------------------|------------------|----------------|-------------------|----------|
| | Nesuit | Qualifier | Office | INDL | - 40101 | | 7a.y = 0 a | | | Allalyst |
| Dissolved Metals by | / MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0958 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:36 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | 0.0079 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | 0.0216 | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 23:01 | EPA 3005A | 64,6020A | ВМ |



12/19/07 13:40

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 Report Date: 01/03/08

SAMPLE RESULTS

Date Collected:

Lab ID: L0718979-03

Client ID: RIZ-9

Date Received: 12/21/07 Sample Location: WALPOLE, MA Field Prep: Field Filtered

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|---------|---------------|-------|--------|-----------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | MCP 600 | 00/7000 serie | es | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0153 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:38 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | ND | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 23:07 | EPA 3005A | 64,6020A | ВМ |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

SAMPLE RESULTS

Lab ID: L0718979-04

Client ID: GHC-6

Sample Location: WALPOLE, MA

ND

Zinc. Dissolved

Matrix: Water

Date Collected: 12/19/07 16:00

Date Received: 12/21/07
Field Prep: Field Filtered

Dilution Analytical Date Date Prep **Factor Prepared** Analyzed Method Method Qualifier **Units RDL Parameter** Result **Analyst** Dissolved Metals by MCP 6000/7000 series Antimony, Dissolved ND 0.0020 64,6020A mg/l 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A BM Arsenic, Dissolved ND 0.0020 4 01/02/08 23:12 EPA 3005A 64,6020A ВМ mg/l 12/30/07 09:30 64,6020A ВМ Barium, Dissolved 0.0459 0.0020 4 01/02/08 23:12 EPA 3005A mg/l 12/30/07 09:30 64,6020A Beryllium, Dissolved ND mg/l 0.0020 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A BM ND 0.0020 4 01/02/08 23:12 EPA 3005A 64,6020A ВМ Cadmium, Dissolved mg/l 12/30/07 09:30 Chromium, Dissolved ND 0.0020 64,6020A ВМ mg/l 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A 64,6020A Lead, Dissolved ND mg/l 0.0020 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A BM Mercury, Dissolved ND mg/l 0.0002 1 12/28/07 16:35 12/31/07 10:40 EPA 7470A 64,7470A DM 64,6020A ND 4 ВМ Nickel, Dissolved 0.0020 01/02/08 23:12 EPA 3005A mg/l 12/30/07 09:30 Selenium, Dissolved ND mg/l 0.004 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A 64,6020A BM 64,6020A Silver, Dissolved ND mg/l 0.0020 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A ВМ Thallium, Dissolved ND 0.0020 4 01/02/08 23:12 EPA 3005A 64,6020A BM mg/l 12/30/07 09:30 Vanadium, Dissolved ND mg/l 0.0020 4 12/30/07 09:30 01/02/08 23:12 EPA 3005A 64,6020A ВМ

mg/l

0.0200

4

12/30/07 09:30



64,6020A

ВМ

01/02/08 23:12 EPA 3005A

12/20/07 09:00

Field Filtered

12/21/07

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L0718979-05

Client ID: MW-9

Sample Location: WALPOLE, MA Field Prep:

Matrix: Water

| _ | | | | | Dilution | Date | Date | Prep | Analytical | |
|----------------------|----------|-------------|-------|--------|----------|----------------|----------------|-----------|------------|---------|
| Parameter | Result | Qualifier | Units | RDL | Factor | Prepared | Analyzed | Method | Method | Analyst |
| Dissolved Metals by | / MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0070 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | BM |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | вм |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | BM |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:42 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | вм |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | 0.0259 | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 23:18 | EPA 3005A | 64,6020A | ВМ |



12/20/07 10:30

Field Filtered

12/21/07

Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Date Collected:

Date Received:

Field Prep:

Lab ID: L0718979-06

Client ID: RIZ-3

Sample Location: WALPOLE, MA

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|---------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | MCP 600 | 00/7000 seri | es | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0256 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:44 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | ND | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 23:23 | EPA 3005A | 64,6020A | ВМ |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

SAMPLE RESULTS

Lab ID: L0718979-07

Client ID: MW-3

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 12/20/07 11:25

Date Received: 12/21/07
Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| | 1100.00 | | | | | | | | | |
| Dissolved Metals by | / MCP 60 | 00/7000 seri | es | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Barium, Dissolved | 0.0152 | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Chromium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Lead, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:45 | EPA 7470A | 64,7470A | DM |
| Nickel, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Selenium, Dissolved | ND | | mg/l | 0.004 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Silver, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |
| Zinc, Dissolved | ND | | mg/l | 0.0200 | 4 | 12/30/07 09:30 | 01/02/08 23:29 | EPA 3005A | 64,6020A | ВМ |



Project Name: WALPOLE PARK SOUTH Lab Number: L0718979

Project Number: 12700058 **Report Date:** 01/03/08

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | _ | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|------------------------|--------------------|------------|------------|--------------------|------------------|------------------|----------------------|----|
| Dissolved Metals by MC | P 6000/7000 series | for sample | e(s): 01-0 | 07 Batch: | WG307204- | 1 | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 12/28/07 16:35 | 12/31/07 10:25 | 64,7470A | DM |

Prep Information

Digestion Method: EPA 7470A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|-----------------------|----------------------|-----------|-----------|--------------------|------------------|------------------|----------------------|---------|
| Dissolved Metals by M | ICP 6000/7000 series | for sampl | e(s): 01- | 07 Batch: | WG307274-1 | ſ | | |
| Antimony, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Barium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Beryllium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Cadmium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Chromium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Lead, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Nickel, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Selenium, Dissolved | ND | mg/l | 0.001 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Silver, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Thallium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |
| Zinc, Dissolved | ND | mg/l | 0.0050 | 1 | 12/30/07 09:30 | 01/02/08 22:12 | 64,6020A | ВМ |

Prep Information

Digestion Method: EPA 3005A



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L0718979

Report Date: 01/03/08

| Parameter | LCS %Recovery | % | LCSD Recovery | %Recovery Limits | RPD | RPD Limits |
|--|----------------------|---------|------------------|-----------------------|-----|------------|
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s) | : 01-07 | Batch: | WG307204-2 WG307204-3 | | |
| Mercury, Dissolved | 98 | | 98 | 80-120 | 0 | 20 |
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s) | : 01-07 | Batch: | WG307274-2 WG307274-3 | | |
| Antimony, Dissolved | 95 | | 98 | 80-120 | 3 | 20 |
| Arsenic, Dissolved | 101 | | 100 | 80-120 | 1 | 20 |
| Barium, Dissolved | 97 | | 100 | 80-120 | 3 | 20 |
| Beryllium, Dissolved | 89 | | 86 | 80-120 | 3 | 20 |
| Cadmium, Dissolved | 104 | | 105 | 80-120 | 1 | 20 |
| Chromium, Dissolved | 101 | | 104 | 80-120 | 3 | 20 |
| Lead, Dissolved | 102 | | 105 | 80-120 | 3 | 20 |
| Nickel, Dissolved | 102 | | 106 | 80-120 | 4 | 20 |
| Selenium, Dissolved | 100 | | 98 | 80-120 | 2 | 20 |
| Silver, Dissolved | 95 | | 97 | 80-120 | 2 | 20 |
| Thallium, Dissolved | 95 | | 99 | 80-120 | 4 | 20 |
| Vanadium, Dissolved | 98 | | 101 | 80-120 | 3 | 20 |
| Zinc, Dissolved | 102 | | 102 | 80-120 | 0 | 20 |



Lab Number: L0718979

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 Report Date: 01/03/08

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal A Absent

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|------------------------------|--------|-----|-------|------|--------|--|
| L0718979-01A | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-01B | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-01C | Plastic 500ml HNO3 preserved | A | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-TL-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-02A | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-02B | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-02C | Plastic 500ml HNO3 preserved | A | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-TL-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-03A | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-03B | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-03C | Plastic 500ml HNO3 preserved | A | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-BA-6020S,MCP-BE-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SB-6020S,MCP-SB-6020S,MCP-SB-6020S,MCP-SB-6020S,MCP-ZN-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-04A | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-04B | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-04C | Plastic 500ml HNO3 preserved | A | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-BE-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-SE-6020S,MCP-ZN-6020S |
| L0718979-05A | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-05B | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |



Project Name: WALPOLE PARK SOUTH

Lab Number: L0718979 Project Number: 12700058 **Report Date:** 01/03/08

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|------------------------------|--------|-----|-------|------|--------|---|
| L0718979-05C | Plastic 500ml HNO3 preserved | A | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-BE-6020S,MCP-CD-6020S,MCP-CR-6020S,MCP-NI-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-SC-6020S,MCP-SC-6020S,MCP-SC-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-06A | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-06B | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-06C | Plastic 500ml HNO3 preserved | А | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-CD-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-SC-6020S,MCP-SC-6020S,MCP-SC-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-07A | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-07B | Vial HCI preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |
| L0718979-07C | Plastic 500ml HNO3 preserved | А | <2 | 2.2 C | Y | Absent | MCP-7470S,MCP-AG-6020S,MCP-AS-6020S,MCP-AS-6020S,MCP-BA-6020S,MCP-CD-6020S,MCP-CR-6020S,MCP-NI-6020S,MCP-PB-6020S,MCP-SB-6020S,MCP-SE-6020S,MCP-SC-6020S,MCP-SC-6020S,MCP-V-6020S,MCP-ZN-6020S |
| L0718979-07N | Vial HCl preserved | Α | N/A | 2.2 C | Υ | Absent | MCP-8260-04 |

Container Comments

L0718979-07B



Project Name:WALPOLE PARK SOUTHLab Number:L0718979Project Number:12700058Report Date:01/03/08

GLOSSARY

Acronyms

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD- Laboratory Control Sample Duplicate: Refer to LCS.

 MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NI - Not Ignitable.

 Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

ND - Not detected at the reported detection limit for the sample.

RDL - Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

The following data qualifiers have been identified for use under the CT DEP Reasonable Confidence Protocols.

- A Spectra identified as "Aldol Condensation Product".
- B The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- J Estimated value. The analyte was tentatively identified; the quantitation is an estimation. (Tentatively identified compounds only.)

Standard Qualifiers

H - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.

Report Format: Data Usability Report



Project Name:WALPOLE PARK SOUTHLab Number:L0718979Project Number:12700058Report Date:01/03/08

REFERENCES

Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). May 2004.

Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Woods Hole Labs performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Woods Hole Labs be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



| PLEASE ANSWER QUESTIONS ABOVE! IS YOUR PROJECT MA MCP or CT RCP? VNO:01-01 (rev. 30-JUL-07) | CHAIN OF CUSTODY ALPHA MESTBORO MA TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TEL: 508-892-9300 TOMMA OF OI Address: |
|---|--|
| Container Type Preservative Relinquished By: Date/Time Date/Jime Date/Jime | CHAIN OF CUSTODY Project Information Project Information Project Information Project Information Project Information Project Mame: Walfele Rock South Project Name: Walfele Rock MA Project Manager: Ray Townson ALPHA Quote #: 200 i ALPHA Quote |
| PX CB Received By: 2014 Sumal | Date Rec'd in La Report Informa Gran ANALYSIS |
| Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguites are resolved. All samples submitted are subject to Alpha's Payment Terms. See reverse side. | b: 12/2/ ALPHA Job #: \(\Lambda \) A \(\text{PP4} \) ASPMITION Are MCP Analytical Methods Required? Are CT RCP (Reasonable Confidence Protocols) Required on Preservation Sample Specific Comments |



ANALYTICAL REPORT

Lab Number: L0806023

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: lan Cannan

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Report Date: 05/05/08

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



L0806023

Lab Number:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 **Report Date:** 05/05/08

| Alpha Sample ID | Client ID | Sample Location |
|-----------------|-------------------|-----------------|
| L0806023-01 | RIZ-10-042808 | WALPOLE, MA |
| L0806023-02 | RIZ-9-042808 | WALPOLE, MA |
| L0806023-03 | MW-3-042808 | WALPOLE, MA |
| L0806023-04 | RIZ-8-042808 | WALPOLE, MA |
| L0806023-05 | RIZ-8S-042808 | WALPOLE, MA |
| L0806023-06 | GHC-6-042808 | WALPOLE, MA |
| L0806023-07 | MW-2-042808 | WALPOLE, MA |
| L0806023-08 | TRIP BLANK-042808 | WALPOLE, MA |

Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| A | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | YES |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



Project Name:WALPOLE PARK SOUTHLab Number:L0806023Project Number:12700058-003Report Date:05/05/08

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

MCP Related Narratives:

Metals

L0806023-01 through -07 were diluted for the analysis of all analytes by method 6020A due to high concentrations of target and non-target analytes.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 05/05/08

ORGANICS



VOLATILES



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-01
 Date Collected:
 04/28/08 09:10

 Client ID:
 RIZ-10-042808
 Date Received:
 04/29/08

Sample Location: WALPOLE, MA Field Prep: Not Specified

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 10:04

Analyst: MM

| Volatile Organics by GC/MS 524.2 ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chioroform ND ug/l 0.50 1 Chiorotorm ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 L2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tichloroflouromethane ND ug/l 0.50 1 Tichloroethane ND ug/l 0.50 1 1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromoform ND ug/l 0.50 | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------------------------|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.5 | Volatile Organics by GC/MS 524.2 | | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorothane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 <td>1,1-Dichloroethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroftuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 dis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ehym-Xylene ND ug/l 0.50 1 | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0,73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Stromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans- | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Interpretation of the companies of the com | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene 0.73 ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | 0.73 | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| · | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-01
 Date Collected:
 04/28/08 09:10

 Client ID:
 RIZ-10-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Tentatively Identified Compounds | | | | |
|----------------------------------|------|---|------|----|
| Unknown Alkene | 2.6 | J | ug/l | 1 |
| Unknown Hydrocarbon | 0.90 | J | ug/l | 1 |
| Acetone | 0.78 | J | ug/l | 11 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 09:10

Client ID:RIZ-10-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| | | Acceptance | | |
|------------------------|------------|------------|----------|--|
| Surrogate | % Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 112 | | 80-120 | |
| 4-Bromofluorobenzene | 88 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-02
 Date Collected:
 04/28/08 10:12

 Client ID:
 RIZ-9-042808
 Date Received:
 04/29/08

Sample Location: WALPOLE, MA Field Prep: Not Specified

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 10:38

Analyst: MM

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-02
 Date Collected:
 04/28/08 10:12

 Client ID:
 RIZ-9-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Not Specified

| Volatile Organics by GC/MS 524.2 1,3-Dichlorobenzene ND | Dilution Factor | RDL | Units | Qualifier | Result | Parameter |
|--|-----------------|------|-------|-----------|--------|----------------------------------|
| 1,4-Dichlorobenzene ND ug/l 0.50 Styrene ND ug/l 0.50 o-Xylene ND ug/l 0.50 1,1-Dichloropropene ND ug/l 0.50 1,2-Dichloropropane ND ug/l 0.50 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 betwicklichtene ND ug/l 0.50 betwicklichtene ND ug/l 0.50 lsopropylbenzene ND ug/l 0.50 lsopropylbenzene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene | | | | | | Volatile Organics by GC/MS 524.2 |
| Styrene ND ug/l 0.50 o-Xylene ND ug/l 0.50 1,1-Dichloropropene ND ug/l 0.50 2,2-Dichloropropane ND ug/l 0.50 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 Ibsopropylbenzene ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene <td>1</td> <td>0.50</td> <td>ug/l</td> <td></td> <td>ND</td> <td>1,3-Dichlorobenzene</td> | 1 | 0.50 | ug/l | | ND | 1,3-Dichlorobenzene |
| o-Xylene ND ug/l 0.50 1,1-Dichloropropene ND ug/l 0.50 2,2-Dichloropropane ND ug/l 0.50 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylteluzene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tet-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trich | 1 | 0.50 | ug/l | | ND | 1,4-Dichlorobenzene |
| 1,1-Dichloropropene ND ug/l 0.50 2,2-Dichloropropane ND ug/l 0.50 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2 | 1 | 0.50 | ug/l | | ND | Styrene |
| 2,2-Dichloropropane ND ug/l 0.50 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 < | 1 | 0.50 | ug/l | | ND | o-Xylene |
| 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 P-Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 Brom | 1 | 0.50 | ug/l | | ND | 1,1-Dichloropropene |
| 1,2,3-Trichloropropane ND ug/l 0.50 Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 Dibromomethane | 1 | 0.50 | ug/l | | ND | 2,2-Dichloropropane |
| Bromochloromethane ND ug/l 0.50 n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 P-Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibrom | 1 | 0.50 | ug/l | | ND | 1,1,1,2-Tetrachloroethane |
| n-Butylbenzene ND ug/l 0.50 Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 P-Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 N-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromo | 1 | 0.50 | ug/l | | ND | 1,2,3-Trichloropropane |
| Dichlorodifluoromethane ND ug/l 0.50 Hexachlorobutadiene ND ug/l 0.50 Isopropylbenzene ND ug/l 0.50 p-Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Bromochloromethane |
| Hexachlorobutadiene ND | 1 | 0.50 | ug/l | | ND | n-Butylbenzene |
| Isopropylbenzene ND | 1 | 0.50 | ug/l | | ND | Dichlorodifluoromethane |
| p-Isopropyltoluene ND ug/l 0.50 Naphthalene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Dibromobenzene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Hexachlorobutadiene |
| Naphthalene ND ug/l 0.50 n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Isopropylbenzene |
| n-Propylbenzene ND ug/l 0.50 sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | p-Isopropyltoluene |
| sec-Butylbenzene ND ug/l 0.50 tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Naphthalene |
| tert-Butylbenzene ND ug/l 0.50 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 pibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | n-Propylbenzene |
| 1,2,3-Trichlorobenzene ND ug/l 0.50 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | sec-Butylbenzene |
| 1,2,4-Trichlorobenzene ND ug/l 0.50 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | tert-Butylbenzene |
| 1,2,4-Trimethylbenzene ND ug/l 0.50 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | 1,2,3-Trichlorobenzene |
| 1,3,5-Trimethylbenzene ND ug/l 0.50 Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | 1,2,4-Trichlorobenzene |
| Bromobenzene ND ug/l 0.50 o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | 1,2,4-Trimethylbenzene |
| o-Chlorotoluene ND ug/l 0.50 p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | 1,3,5-Trimethylbenzene |
| p-Chlorotoluene ND ug/l 0.50 Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Bromobenzene |
| Dibromomethane ND ug/l 0.50 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | o-Chlorotoluene |
| 1,2-Dibromoethane ND ug/l 0.50 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | p-Chlorotoluene |
| 1,2-Dibromo-3-chloropropane ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Dibromomethane |
| | 1 | 0.50 | ug/l | | ND | 1,2-Dibromoethane |
| 1.3 Dichloropropage | 1 | 0.50 | ug/l | | ND | 1,2-Dibromo-3-chloropropane |
| 1,3-Dictilioroproparie ND ug/i 0.30 | 1 | 0.50 | ug/l | | ND | 1,3-Dichloropropane |
| Methyl tert butyl ether ND ug/l 0.50 | 1 | 0.50 | ug/l | | ND | Methyl tert butyl ether |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-02 Date Collected: 04/28/08 10:12

Client ID:RIZ-9-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 116 | | 80-120 | |
| 4-Bromofluorobenzene | 87 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-03 Date Collected: 04/28/08 11:20

Client ID:MW-3-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 11:10

Analyst: MM

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-03
 Date Collected:
 04/28/08 11:20

 Client ID:
 MW-3-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Not Specified

| Volatile Organics by GC/MS 524.2 1,3-Dichlorobenzene ND ug/l 0.50 1 1,4-Dichlorobenzene ND ug/l 0.50 1 Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 I-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloropethane ND ug/l 0.50 1 1,1,1,2-Tetrachloropethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropyltoluene ND ug/l 0.50 1 Hexachlorobutaen | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------------------------|--------|-----------|-------|------|-----------------|
| 1,4-Dichlorobenzene ND ug/l 0.50 1 Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 1,1-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloropthane ND ug/l 0.50 1 1,2,3-Trichloropthane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 NButylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadilene ND ug/l 0.50 1 Bopropylbenzene ND ug/l 0.50 1 p-Isopropylbeluzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 | Volatile Organics by GC/MS 524.2 | | | | | |
| Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropyllouzene ND ug/l 0.50 1 Isopropyllouzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 | 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylibenzene ND ug/l 0.50 1 Isopropylibenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 | 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodiffluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 <t< td=""><td>Styrene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Styrene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 <t< td=""><td>o-Xylene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 <td< td=""><td>1,1-Dichloropropene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane ND | 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Let-Butylbenzene ND ug/l 0.50 1 Let-Butylbe | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 Letr-Butylbenzene ND ug/l 0.50 1 Letr-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 O-Chlorotoluene ND ug/l 0.50 1 | 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1< | Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene ND ug/l 0.50 1 | n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Sopropylbenzene ND | Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 | Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 | Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Naphthalene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 1 o-Chlorotoluene ND ug/l 0.50 1 | n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Bromobenzene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether ND ug/l 0.50 1 | 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| | Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-03 Date Collected: 04/28/08 11:20

Client ID:MW-3-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 112 | | 80-120 | |
| 4-Bromofluorobenzene | 90 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 12:15

Client ID:RIZ-8-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 11:43

Analyst: MM

| Volatile Organics by GC/MS 524.2 ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 L2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tichlorofobanzene ND ug/l 0.50 1 Tichloroethane ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------------------------|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 L1,2-Dichloropthane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 tetras-1,3-Dichloropropene ND ug/l 0.50 1 gis-1,3-Dichloropropene ND ug/l 0.50 | Volatile Organics by GC/MS 524.2 | | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 <t< td=""><td>Methylene chloride</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorothane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylsenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroftuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 sis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 < | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <t< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Stromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1, | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 <t< td=""><td>1,2-Dichloroethane</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Interpretation of the control o | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| · | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-04
 Date Collected:
 04/28/08 12:15

 Client ID:
 RIZ-8-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Field Filtered

| Volatile Organics by GC/MS 524.2 1,3-Dichlorobenzene ND ug/l 0.50 1 1,4-Dichlorobenzene ND ug/l 0.50 1 Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 I-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloropethane ND ug/l 0.50 1 1,1,1,2-Tetrachloropethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropyltoluene ND ug/l 0.50 1 Hexachlorobutaen | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------------------------|--------|-----------|-------|------|-----------------|
| 1,4-Dichlorobenzene ND ug/l 0.50 1 Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 1,1-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloropthane ND ug/l 0.50 1 1,2,3-Trichloropthane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 NButylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadilene ND ug/l 0.50 1 Bopropylbenzene ND ug/l 0.50 1 p-Isopropylbeluzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 | Volatile Organics by GC/MS 524.2 | | | | | |
| Styrene ND ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropyllouzene ND ug/l 0.50 1 Isopropyllouzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 | 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylibenzene ND ug/l 0.50 1 Isopropylibenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 | 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodiffluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 <t< td=""><td>Styrene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Styrene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Ibopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 <t< td=""><td>o-Xylene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 1 tetr-Butylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 <td< td=""><td>1,1-Dichloropropene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane ND | 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Let-Butylbenzene ND ug/l 0.50 1 Let-Butylbe | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 P-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 Letr-Butylbenzene ND ug/l 0.50 1 Letr-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 O-Chlorotoluene ND ug/l 0.50 1 | 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1< | Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene ND ug/l 0.50 1 | n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Sopropylbenzene ND | Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 | Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 | Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Naphthalene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 1 o-Chlorotoluene ND ug/l 0.50 1 | n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Bromobenzene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether ND ug/l 0.50 1 | 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| | Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-04 Date Collected: 04/28/08 12:15

Client ID:RIZ-8-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 114 | | 80-120 | |
| 4-Bromofluorobenzene | 91 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

SAMPLE RESULTS

 Lab ID:
 L0806023-05
 Date Collected:
 04/28/08 12:30

 Client ID:
 RIZ-8S-042808
 Date Received:
 04/29/08

Client ID:RIZ-8S-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 05/01/08 12:17

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 11 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-05
 Date Collected:
 04/28/08 12:30

 Client ID:
 RIZ-8S-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 11 |
| | | | | | |

| Tentatively Identified Compounds | | | | |
|----------------------------------|------|---|------|----|
| Unknown Alkane | 0.62 | J | ug/l | 11 |
| Unknown Alkane | 2.7 | J | ug/l | 1 |
| Unknown Hydrocarbon | 4.3 | J | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-05 Date Collected: 04/28/08 12:30

Client ID: RIZ-8S-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Field Filtered

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

 Tentatively Identified Compounds

 Unknown Hydrocarbon
 0.64
 J
 ug/l
 1

 Pentane
 1.2
 J
 ug/l
 1

Surrogate % Recovery Qualifier Criteria

1,2-Dichlorobenzene-d4 116 80-120

4-Bromofluorobenzene 87 80-120



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-06 Date Collected: 04/28/08 14:36

Client ID:GHC-6-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 12:50

| Volatile Organics by GC/MS 524.2 ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 L2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tichlorofobanzene ND ug/l 0.50 1 Tichloroethane ND ug/l 0.50 1 Tichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Intans-1,3-Dichloropropene ND ug/l 0.50 | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------------------------|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Ly-Eichloroethane ND ug/l 0.50 1 Eromodichloromethane ND ug/l 0.50 1 Eromodichloropropene ND ug/l 0.50 1 Bromodichm ND ug/l 0.50 1 | Volatile Organics by GC/MS 524.2 | | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 <t< td=""><td>Methylene chloride</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorothane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylsenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroftuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 sis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 < | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <t< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Stromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1, | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 <t< td=""><td>1,2-Dichloroethane</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></t<> | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Interpretation of the control o | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| · | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-06
 Date Collected:
 04/28/08 14:36

 Client ID:
 GHC-6-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-06 Date Collected: 04/28/08 14:36

Client ID:GHC-6-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 117 | | 80-120 | |
| 4-Bromofluorobenzene | 88 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 14:52

Client ID:MW-2-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 13:23

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | 2.2 | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

 Lab ID:
 L0806023-07
 Date Collected:
 04/28/08 14:52

 Client ID:
 MW-2-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lo806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-07 Date Collected: 04/28/08 14:52

Client ID:MW-2-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Field Filtered

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 118 | | 80-120 | |
| 4-Bromofluorobenzene | 87 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-08 Date Collected: 04/23/08 18:00

Client ID:TRIP BLANK-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Matrix: Water
Anaytical Method: 16,524.2
Analytical Date: 05/01/08 09:31

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|----------------------------------|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS 524.2 | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

SAMPLE RESULTS

 Lab ID:
 L0806023-08
 Date Collected:
 04/23/08 18:00

 Client ID:
 TRIP BLANK-042808
 Date Received:
 04/29/08

 Sample Location:
 WALPOLE, MA
 Field Prep:
 Not Specified

Parameter Result Qualifier Units **RDL Dilution Factor** Volatile Organics by GC/MS 524.2 1,3-Dichlorobenzene ND 0.50 ug/l 1 ND 1 1,4-Dichlorobenzene ug/l 0.50 ND Styrene ug/l 0.50 1 o-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND 0.50 ug/l 1 ND 2,2-Dichloropropane ug/l 0.50 1 ND 1 1,1,1,2-Tetrachloroethane ug/l 0.50 ND 1 1,2,3-Trichloropropane ug/l 0.50 ND Bromochloromethane ug/l 0.50 1 ND n-Butylbenzene ug/l 0.50 1 Dichlorodifluoromethane ND 0.50 ug/l 1 Hexachlorobutadiene ND ug/l 0.50 1 ND 1 Isopropylbenzene ug/l 0.50 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND 0.50 ug/l 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 ND 1,2,4-Trichlorobenzene ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 ND 1,3,5-Trimethylbenzene ug/l 0.50 1 ND Bromobenzene ug/l 0.50 1 ND 1 o-Chlorotoluene ug/l 0.50 ND 1 p-Chlorotoluene ug/l 0.50 Dibromomethane ND ug/l 0.50 1 ND 0.50 1 1,2-Dibromoethane ug/l 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 ND Methyl tert butyl ether ug/l 0.50 1

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-08 Date Collected: 04/23/08 18:00

Client ID:TRIP BLANK-042808Date Received:04/29/08Sample Location:WALPOLE, MAField Prep:Not Specified

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS 524.2

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 113 | | 80-120 | |
| 4-Bromofluorobenzene | 95 | | 80-120 | |



Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

| Parameter | Result | Qualifier | Units | RDL |
|---------------------------------|-----------------|------------|--------|------------|
| olatile Organics by GC/MS 524.2 | 2 for sample(s) | : 01,04-08 | Batch: | WG319665-8 |
| Methylene chloride | ND | | ug/l | 0.50 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 |
| Chloroform | ND | | ug/l | 0.50 |
| Carbon tetrachloride | ND | | ug/l | 0.50 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 |
| Dibromochloromethane | ND | | ug/l | 0.50 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 |
| Tetrachloroethene | ND | | ug/l | 0.50 |
| Chlorobenzene | ND | | ug/l | 0.50 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 |
| Bromodichloromethane | ND | | ug/l | 0.50 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 |
| Bromoform | ND | | ug/l | 0.50 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 |
| Benzene | ND | | ug/l | 0.50 |
| Toluene | ND | | ug/l | 0.50 |
| Ethylbenzene | ND | | ug/l | 0.50 |
| p/m-Xylene | ND | | ug/l | 0.50 |
| Chloromethane | ND | | ug/l | 0.50 |
| Bromomethane | ND | | ug/l | 0.50 |
| Vinyl chloride | ND | | ug/l | 0.50 |
| Chloroethane | ND | | ug/l | 0.50 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 |
| Trichloroethene | ND | | ug/l | 0.50 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 |



Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

| arameter | Result | Qualifier | Units | RDL |
|---------------------------------|---------------|------------|--------|------------|
| olatile Organics by GC/MS 524.2 | for sample(s) | : 01,04-08 | Batch: | WG319665-8 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 |
| Styrene | ND | | ug/l | 0.50 |
| o-Xylene | ND | | ug/l | 0.50 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 |
| Bromochloromethane | ND | | ug/l | 0.50 |
| n-Butylbenzene | ND | | ug/l | 0.50 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 |
| Isopropylbenzene | ND | | ug/l | 0.50 |
| p-lsopropyltoluene | ND | | ug/l | 0.50 |
| Naphthalene | ND | | ug/l | 0.50 |
| n-Propylbenzene | ND | | ug/l | 0.50 |
| sec-Butylbenzene | ND | | ug/l | 0.50 |
| tert-Butylbenzene | ND | | ug/l | 0.50 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 |
| Bromobenzene | ND | | ug/l | 0.50 |
| o-Chlorotoluene | ND | | ug/l | 0.50 |
| p-Chlorotoluene | ND | | ug/l | 0.50 |
| Dibromomethane | ND | | ug/l | 0.50 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 |



Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

Analyst: MM

Parameter Result Qualifier Units RDL

Volatile Organics by GC/MS 524.2 for sample(s): 01,04-08 Batch: WG319665-8

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | Acceptance | | | |
|------------------------|------------|-----------|----------|--|
| Surrogate | %Recovery | Qualifier | Criteria | |
| | | | | |
| 1,2-Dichlorobenzene-d4 | 114 | | 80-120 | |
| 4-Bromofluorobenzene | 87 | | 80-120 | |



Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

| Result | Qualifier | Unit | s RDL | |
|----------------|--|--|--|---|
| for sample(s): | 02-03 | Batch: | WG320296-4 | |
| ND | | ug/l | 0.50 | |
| ND | | | | |
| ND | | | | |
| ND | | | | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| ND | | ug/l | 0.50 | |
| | ND ND ND ND ND ND ND ND ND ND ND ND ND N | ND ND ND ND ND ND ND ND ND ND ND ND ND N | ND ug/l ND | ND ug/l 0.50 ND ug/l |



Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

| Result | Qualifier | Unit | s RDL |
|---------------|--|--|--|
| for sample(s) | : 02-03 | Batch: | WG320296-4 |
| ND | | ug/l | 0.50 |
| | For sample(s) ND ND ND ND ND ND ND ND ND N | ND ND ND ND ND ND ND ND ND ND ND ND ND N | ND ug/l ND |



Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 05/01/08 08:25

Analyst: MM

Parameter Result Qualifier Units RDL

Volatile Organics by GC/MS 524.2 for sample(s): 02-03 Batch: WG320296-4

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | | Acceptance | | | |
|------------------------|-----------|------------|----------|--|--|
| Surrogate | %Recovery | Qualifier | Criteria | | |
| | | | | | |
| 1,2-Dichlorobenzene-d4 | 114 | | 80-120 | | |
| 4-Bromofluorobenzene | 87 | | 80-120 | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 Lab Number:

L0806023

| Methylene chloride 1,1-Dichloroethane Chloroform Carbon tetrachloride 1,2-Dichloropropane Dibromochloromethane | 99 107 106 | 08 Batch: WG319665- - - | -7 70-130 | <u>-</u> | |
|--|------------------|-------------------------------|------------------|--------------|--|
| 1,1-Dichloroethane Chloroform Carbon tetrachloride 1,2-Dichloropropane | 107 106 | - | 70-130 | - | |
| Chloroform Carbon tetrachloride 1,2-Dichloropropane | 106 | - | | | |
| Carbon tetrachloride 1,2-Dichloropropane | | | 70-130 | - | |
| 1,2-Dichloropropane | | - | 70-130 | - | |
| | 103 | - | 70-130 | - | |
| Dibromochloromethane | 106 | - | 70-130 | - | |
| | 96 | - | 70-130 | - | |
| 1,1,2-Trichloroethane | 98 | - | 70-130 | - | |
| Tetrachloroethene | 112 | - | 70-130 | - | |
| Chlorobenzene | 108 | - | 70-130 | - | |
| Trichlorofluoromethane | 93 | - | 70-130 | - | |
| 1,2-Dichloroethane | 109 | - | 70-130 | - | |
| 1,1,1-Trichloroethane | 108 | - | 70-130 | - | |
| Bromodichloromethane | 102 | - | 70-130 | - | |
| trans-1,3-Dichloropropene | 86 | - | 70-130 | - | |
| cis-1,3-Dichloropropene | 90 | - | 70-130 | - | |
| Bromoform | 90 | - | 70-130 | - | |
| 1,1,2,2-Tetrachloroethane | 90 | - | 70-130 | - | |
| Benzene | 111 | - | 70-130 | - | |
| Toluene | 116 | - | 70-130 | | |
| Ethylbenzene | 110 | | 70-130 | - | |
| p/m-Xylene | 110 | - | 70-130 | - | |

WALPOLE PARK SOUTH

Project Number: 12700058-003

Project Name:

Lab Number:

L0806023

Report Date:

05/05/08

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---------------------------------|-------------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by GC/MS 524.2 | Associated sample(s): 01,04-0 | 8 Batch: WG319665 | 5-7 | | |
| Chloromethane | 101 | - | 70-130 | - | |
| Bromomethane | 103 | - | 70-130 | - | |
| Vinyl chloride | 107 | - | 70-130 | - | |
| Chloroethane | 106 | - | 70-130 | - | |
| 1,1-Dichloroethene | 100 | - | 70-130 | - | |
| trans-1,2-Dichloroethene | 103 | - | 70-130 | - | |
| cis-1,2-Dichloroethene | 111 | - | 70-130 | - | |
| Trichloroethene | 101 | - | 70-130 | - | |
| 1,2-Dichlorobenzene | 95 | - | 70-130 | - | |
| 1,3-Dichlorobenzene | 99 | - | 70-130 | - | |
| 1,4-Dichlorobenzene | 95 | - | 70-130 | - | |
| Styrene | 108 | - | 70-130 | - | |
| o-Xylene | 109 | - | 70-130 | - | |
| 1,1-Dichloropropene | 107 | - | 70-130 | - | |
| 2,2-Dichloropropane | 93 | - | 70-130 | - | |
| 1,1,1,2-Tetrachloroethane | 101 | - | 70-130 | - | |
| 1,2,3-Trichloropropane | 91 | - | 70-130 | - | |
| Bromochloromethane | 106 | - | 70-130 | - | |
| n-Butylbenzene | 102 | - | 70-130 | - | |
| Dichlorodifluoromethane | 106 | - | 70-130 | - | |
| Hexachlorobutadiene | 98 | - | 70-130 | - | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| nrameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---------------------------------|--------------------------------|-------------------|---------------------|-----|------------|
| platile Organics by GC/MS 524.2 | Associated sample(s): 01,04-08 | Batch: WG319665 | 5-7 | | |
| Isopropylbenzene | 103 | - | 70-130 | - | |
| p-Isopropyltoluene | 103 | - | 70-130 | - | |
| Naphthalene | 82 | - | 70-130 | - | |
| n-Propylbenzene | 113 | - | 70-130 | - | |
| sec-Butylbenzene | 109 | - | 70-130 | - | |
| tert-Butylbenzene | 112 | - | 70-130 | - | |
| 1,2,3-Trichlorobenzene | 87 | - | 70-130 | - | |
| 1,2,4-Trichlorobenzene | 86 | - | 70-130 | - | |
| 1,2,4-Trimethylbenzene | 106 | - | 70-130 | - | |
| 1,3,5-Trimethylbenzene | 115 | - | 70-130 | - | |
| Bromobenzene | 103 | - | 70-130 | - | |
| o-Chlorotoluene | 110 | - | 70-130 | - | |
| p-Chlorotoluene | 105 | - | 70-130 | - | |
| Dibromomethane | 102 | - | 70-130 | - | |
| 1,2-Dibromoethane | 92 | - | 70-130 | - | |
| 1,2-Dibromo-3-chloropropane | 78 | - | 70-130 | - | |
| 1,3-Dichloropropane | 98 | - | 70-130 | - | |
| Methyl tert butyl ether | 97 | - | 70-130 | - | |

Lab Number: L0806023

Report Date: 05/05/08

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

LCS **LCSD** %Recovery Limits %Recovery %Recovery Parameter

RPD **RPD Limits**

Volatile Organics by GC/MS 524.2 Associated sample(s): 01,04-08 Batch: WG319665-7

| Surrogate | LCS %Recovery Qualifier | LCSD %Recovery Qualifier | Acceptance Criteria | |
|------------------------|----------------------------|-----------------------------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 102 | | 80-120 | |

| Volatile Organics by GC/MS 524.2 Associated | d sample(s): 02 | 2-03 Batch: WG320296-3 | | |
|---|-----------------|------------------------|--------|---|
| Methylene chloride | 99 | - | 70-130 | - |
| 1,1-Dichloroethane | 107 | - | 70-130 | - |
| Chloroform | 106 | - | 70-130 | - |
| Carbon tetrachloride | 103 | - | 70-130 | - |
| 1,2-Dichloropropane | 106 | - | 70-130 | - |
| Dibromochloromethane | 96 | - | 70-130 | - |
| 1,1,2-Trichloroethane | 98 | - | 70-130 | - |
| Tetrachloroethene | 112 | - | 70-130 | - |
| Chlorobenzene | 108 | - | 70-130 | - |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---------------------------------------|------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by GC/MS 524.2 Assoc | iated sample(s): 02-03 | Batch: WG320296-3 | | | |
| Trichlorofluoromethane | 93 | - | 70-130 | - | |
| 1,2-Dichloroethane | 109 | - | 70-130 | - | |
| 1,1,1-Trichloroethane | 108 | - | 70-130 | - | |
| Bromodichloromethane | 102 | - | 70-130 | - | |
| trans-1,3-Dichloropropene | 86 | - | 70-130 | - | |
| cis-1,3-Dichloropropene | 90 | - | 70-130 | - | |
| Bromoform | 90 | - | 70-130 | - | |
| 1,1,2,2-Tetrachloroethane | 90 | - | 70-130 | - | |
| Benzene | 111 | - | 70-130 | - | |
| Toluene | 116 | - | 70-130 | - | |
| Ethylbenzene | 110 | - | 70-130 | - | |
| p/m-Xylene | 113 | - | 70-130 | - | |
| Chloromethane | 101 | - | 70-130 | - | |
| Bromomethane | 103 | - | 70-130 | - | |
| Vinyl chloride | 107 | - | 70-130 | - | |
| Chloroethane | 106 | - | 70-130 | - | |
| 1,1-Dichloroethene | 100 | - | 70-130 | - | |
| trans-1,2-Dichloroethene | 103 | - | 70-130 | - | |
| cis-1,2-Dichloroethene | 111 | - | 70-130 | - | |
| Trichloroethene | 101 | - | 70-130 | - | |
| 1,2-Dichlorobenzene | 95 | - | 70-130 | - | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|-----------------------------------|----------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by GC/MS 524.2 A | ssociated sample(s): 02-03 | Batch: WG320296-3 | | | |
| 1,3-Dichlorobenzene | 99 | - | 70-130 | - | |
| 1,4-Dichlorobenzene | 95 | - | 70-130 | - | |
| Styrene | 108 | - | 70-130 | - | |
| o-Xylene | 109 | - | 70-130 | - | |
| 1,1-Dichloropropene | 107 | - | 70-130 | - | |
| 2,2-Dichloropropane | 93 | - | 70-130 | - | |
| 1,1,1,2-Tetrachloroethane | 101 | - | 70-130 | - | |
| 1,2,3-Trichloropropane | 91 | - | 70-130 | - | |
| Bromochloromethane | 106 | - | 70-130 | - | |
| n-Butylbenzene | 102 | - | 70-130 | - | |
| Dichlorodifluoromethane | 106 | - | 70-130 | - | |
| Hexachlorobutadiene | 98 | - | 70-130 | - | |
| Isopropylbenzene | 103 | - | 70-130 | - | |
| p-Isopropyltoluene | 103 | - | 70-130 | - | |
| Naphthalene | 82 | - | 70-130 | - | |
| n-Propylbenzene | 113 | - | 70-130 | - | |
| sec-Butylbenzene | 109 | - | 70-130 | - | |
| tert-Butylbenzene | 112 | - | 70-130 | - | |
| 1,2,3-Trichlorobenzene | 87 | - | 70-130 | - | |
| 1,2,4-Trichlorobenzene | 86 | - | 70-130 | - | |
| 1,2,4-Trimethylbenzene | 106 | - | 70-130 | - | |

Lab Number:

L0806023

05/05/08

Project Number: 12700058-003

WALPOLE PARK SOUTH

Project Name:

Report Date:

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|-------------------------------------|---------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by GC/MS 524.2 Ass | sociated sample(s): 02-03 | Batch: WG320296-3 | | | |
| 1,3,5-Trimethylbenzene | 115 | - | 70-130 | - | |
| Bromobenzene | 103 | - | 70-130 | - | |
| o-Chlorotoluene | 110 | - | 70-130 | - | |
| p-Chlorotoluene | 105 | - | 70-130 | - | |
| Dibromomethane | 102 | - | 70-130 | - | |
| 1,2-Dibromoethane | 92 | - | 70-130 | - | |
| 1,2-Dibromo-3-chloropropane | 78 | - | 70-130 | - | |
| 1,3-Dichloropropane | 98 | - | 70-130 | - | |
| Methyl tert butyl ether | 97 | - | 70-130 | - | |

| Surrogate | LCS %Recovery Qualifier | LCSD %Recovery Qualifier | Acceptance Criteria | |
|------------------------|----------------------------|-----------------------------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 102 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|----------------------------|------------------|---------------|-----------|-----------------|------------|------------------|--------------------|-----------|------------|
| Volatile Organics by GC/MS | 524.2 Associated | sample(s): 01 | ,04-08 QC | Batch ID: WG3 | 19665-1 QC | Sample: L080 | 5753-01 Cli | ent ID: M | 1S Sample |
| Methylene chloride | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| Chloroform | ND | 4 | 4.4 | 109 | - | - | 70-130 | - | 20 |
| Carbon tetrachloride | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.2 | 105 | - | - | 70-130 | - | 20 |
| Dibromochloromethane | ND | 4 | 3.9 | 97 | - | - | 70-130 | - | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Tetrachloroethene | ND | 4 | 4.7 | 117 | - | - | 70-130 | - | 20 |
| Chlorobenzene | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| Trichlorofluoromethane | ND | 4 | 4.1 | 103 | - | - | 70-130 | - | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.4 | 109 | - | - | 70-130 | - | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| Bromodichloromethane | ND | 4 | 4.2 | 105 | - | - | 70-130 | - | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 3.4 | 85 | - | - | 70-130 | - | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 3.6 | 90 | - | - | 70-130 | - | 20 |
| Bromoform | ND | 4 | 3.8 | 96 | - | - | 70-130 | - | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| Benzene | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| Toluene | ND | 4 | 4.8 | 121 | - | - | 70-130 | - | 20 |
| Ethylbenzene | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| p/m-Xylene | ND | 8 | 9.7 | 121 | - | - | 70-130 | - | 20 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Foun | MSD d %Recovery | Recovery Limits | / RPD | RPD Limits |
|----------------------------|------------------|---------------|-----------|-----------------|----------|--------------------|--------------------|--------------|------------|
| Volatile Organics by GC/MS | 524.2 Associated | sample(s): 01 | ,04-08 QC | Batch ID: WG: | 319665-1 | QC Sample: L080 | 5753-01 (| Client ID: M | 1S Sample |
| Chloromethane | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |
| Bromomethane | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Vinyl chloride | ND | 4 | 4.7 | 118 | - | - | 70-130 | - | 20 |
| Chloroethane | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.2 | 104 | - | - | 70-130 | - | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| Trichloroethene | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 4.1 | 104 | - | - | 70-130 | - | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Styrene | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| o-Xylene | ND | 4 | 4.7 | 118 | - | - | 70-130 | - | 20 |
| 1,1-Dichloropropene | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| 2,2-Dichloropropane | ND | 4 | 3.8 | 95 | - | - | 70-130 | - | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 4.3 | 109 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| Bromochloromethane | ND | 4 | 4.2 | 105 | - | - | 70-130 | - | 20 |
| n-Butylbenzene | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| Dichlorodifluoromethane | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| Hexachlorobutadiene | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|-----------------------------|------------------|---------------|-----------|-----------------|-------------|------------------|--------------------|------------|------------|
| olatile Organics by GC/MS | 524.2 Associated | sample(s): 01 | ,04-08 QC | Batch ID: WG | 319665-1 Q0 | C Sample: L080 | 5753-01 CI | ient ID: M | IS Sample |
| Isopropylbenzene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| p-Isopropyltoluene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| Naphthalene | ND | 4 | 3.6 | 89 | - | - | 70-130 | - | 20 |
| n-Propylbenzene | ND | 4 | 4.9 | 123 | - | - | 70-130 | - | 20 |
| sec-Butylbenzene | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| tert-Butylbenzene | ND | 4 | 4.9 | 123 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 4.9 | 122 | - | - | 70-130 | - | 20 |
| Bromobenzene | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| o-Chlorotoluene | ND | 4 | 4.7 | 118 | - | - | 70-130 | - | 20 |
| p-Chlorotoluene | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| Dibromomethane | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |
| 1,2-Dibromoethane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| 1,3-Dichloropropane | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| Methyl tert butyl ether | ND | 4 | 3.6 | 90 | - | - | 70-130 | - | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| MS MSD Recov | ery |
|--------------|-----|
|--------------|-----|

Parameter Native Sample MS Added MS Found %Recovery MSD Found %Recovery Limits RPD RPD Limits

Volatile Organics by GC/MS 524.2 Associated sample(s): 01,04-08 QC Batch ID: WG319665-1 QC Sample: L0805753-01 Client ID: MS Sample

| | MS | ; | M | SD | Acceptance | |
|------------------------|------------|-----------|------------|-----------|------------|--|
| Surrogate | % Recovery | Qualifier | % Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 101 | | | | 80-120 | |
| 4-Bromofluorobenzene | 104 | | | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| Parameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Four | MSD nd %Recovery | Recovery Limits | RPD | RPD Limits |
|----------------------------|------------------|----------------|------------|-----------------|----------|---------------------|--------------------|--------|------------|
| Volatile Organics by GC/MS | 524.2 Associated | sample(s): 02- | 03 QC Bate | ch ID: WG3202 | 296-1 QC | Sample: L080602 | 3-02 Client IE | : RIZ- | 9-042808 |
| Methylene chloride | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.6 | 116 | - | - | 70-130 | - | 20 |
| Chloroform | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| Carbon tetrachloride | ND | 4 | 4.7 | 119 | - | - | 70-130 | - | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| Dibromochloromethane | ND | 4 | 4.1 | 103 | - | - | 70-130 | - | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |
| Tetrachloroethene | ND | 4 | 5.0 | 124 | - | - | 70-130 | - | 20 |
| Chlorobenzene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| Trichlorofluoromethane | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 4.8 | 119 | - | - | 70-130 | - | 20 |
| Bromodichloromethane | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 3.6 | 91 | - | - | 70-130 | - | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| Bromoform | ND | 4 | 3.8 | 95 | - | - | 70-130 | - | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 3.8 | 96 | - | - | 70-130 | - | 20 |
| Benzene | ND | 4 | 5.0 | 125 | - | - | 70-130 | - | 20 |
| Toluene | ND | 4 | 5.2 | 130 | - | - | 70-130 | - | 20 |
| Ethylbenzene | ND | 4 | 4.7 | 117 | - | - | 70-130 | - | 20 |
| p/m-Xylene | ND | 8 | 9.6 | 120 | - | - | 70-130 | - | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| Parameter | Native Sample | MS Added M | IS Found | MS %Recovery N | /ISD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|----------------------------|------------------|------------------|-----------|-------------------|------------|------------------|--------------------|----------|------------|
| Volatile Organics by GC/MS | 524.2 Associated | sample(s): 02-03 | 3 QC Batc | h ID: WG32029 | 6-1 QC Sa | mple: L080602 | 3-02 Client I | D: RIZ-9 | 9-042808 |
| Chloromethane | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| Bromomethane | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| Vinyl chloride | ND | 4 | 4.9 | 122 | - | - | 70-130 | - | 20 |
| Chloroethane | ND | 4 | 4.9 | 123 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| Trichloroethene | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 3.9 | 97 | - | - | 70-130 | - | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 3.9 | 97 | - | - | 70-130 | - | 20 |
| Styrene | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| o-Xylene | ND | 4 | 4.6 | 116 | - | - | 70-130 | - | 20 |
| 1,1-Dichloropropene | ND | 4 | 4.7 | 118 | - | - | 70-130 | - | 20 |
| 2,2-Dichloropropane | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| Bromochloromethane | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| n-Butylbenzene | ND | 4 | 4.1 | 103 | - | - | 70-130 | - | 20 |
| Dichlorodifluoromethane | ND | 4 | 4.9 | 122 | - | - | 70-130 | - | 20 |
| Hexachlorobutadiene | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| arameter | Native Sample | MS Added M | IS Found % | MS %Recovery M | /ISD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|-----------------------------|--------------------|------------------|------------|-------------------|------------|------------------|--------------------|---------|------------|
| olatile Organics by GC/MS | S 524.2 Associated | sample(s): 02-03 | 3 QC Batch | n ID: WG32029 | 6-1 QC Sa | mple: L080602 | 3-02 Client I | D: RIZ- | 9-042808 |
| Isopropylbenzene | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| p-Isopropyltoluene | ND | 4 | 4.3 | 109 | - | - | 70-130 | - | 20 |
| Naphthalene | ND | 4 | 3.3 | 83 | - | - | 70-130 | - | 20 |
| n-Propylbenzene | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| sec-Butylbenzene | ND | 4 | 4.7 | 117 | - | <u>-</u> | 70-130 | - | 20 |
| tert-Butylbenzene | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.6 | 90 | - | - | 70-130 | - | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.5 | 87 | - | - | 70-130 | - | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 4.8 | 119 | - | - | 70-130 | - | 20 |
| Bromobenzene | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| o-Chlorotoluene | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| p-Chlorotoluene | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| Dibromomethane | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| 1,2-Dibromoethane | ND | 4 | 3.8 | 96 | - | - | 70-130 | - | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 3.2 | 81 | - | - | 70-130 | - | 20 |
| 1,3-Dichloropropane | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |
| Methyl tert butyl ether | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| MS MSD | Recovery |
|--------|----------|
|--------|----------|

Parameter Native Sample MS Added MS Found %Recovery MSD Found %Recovery Limits RPD RPD Limits

Volatile Organics by GC/MS 524.2 Associated sample(s): 02-03 QC Batch ID: WG320296-1 QC Sample: L0806023-02 Client ID: RIZ-9-042808

| | MS | M | SD | Acceptance | |
|------------------------|----------------|--------------------|-----------|------------|--|
| Surrogate | % Recovery Qua | alifier % Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 98 | | | 80-120 | |
| 4-Bromofluorobenzene | 103 | | | 80-120 | |



Lab Duplicate Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| rameter | Native S | ample Duplicate S | Sample Units | RPD | RPD Limits |
|---------------------------------|--------------------------------|----------------------|----------------|-----------------------|---------------|
| platile Organics by GC/MS 524.2 | Associated sample(s): 01,04-08 | QC Batch ID: WG31966 | 5-2 QC Sample: | L0805753-02 Client IE | D: DUP Sample |
| Methylene chloride | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethane | ND | ND | ug/l | NC | 20 |
| Chloroform | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene | ND | ND | ug/l | NC | 20 |
| Chlorobenzene | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| Bromoform | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| Benzene | ND | ND | ug/l | NC | 20 |
| Toluene | ND | ND | ug/l | NC | 20 |



Lab Duplicate Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| arameter | N | lative Sar | nple Duplicate S | ample Units | s RPD | RPD Limits |
|---------------------------------|-------------------------|------------|-----------------------|----------------|--------------------|----------------|
| olatile Organics by GC/MS 524.2 | Associated sample(s): 0 | 01,04-08 | QC Batch ID: WG319665 | 5-2 QC Sample: | L0805753-02 Client | ID: DUP Sample |
| Ethylbenzene | | ND | ND | ug/l | NC | 20 |
| p/m-Xylene | | ND | ND | ug/l | NC | 20 |
| Chloromethane | | ND | ND | ug/l | NC | 20 |
| Bromomethane | | ND | ND | ug/l | NC | 20 |
| Vinyl chloride | | ND | ND | ug/l | NC | 20 |
| Chloroethane | | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethene | | ND | ND | ug/l | NC | 20 |
| trans-1,2-Dichloroethene | | ND | ND | ug/l | NC | 20 |
| cis-1,2-Dichloroethene | | ND | ND | ug/l | NC | 20 |
| Trichloroethene | | ND | ND | ug/l | NC | 20 |
| 1,2-Dichlorobenzene | | ND | ND | ug/l | NC | 20 |
| 1,3-Dichlorobenzene | | ND | ND | ug/l | NC | 20 |
| 1,4-Dichlorobenzene | | ND | ND | ug/l | NC | 20 |
| Styrene | | ND | ND | ug/l | NC | 20 |
| o-Xylene | | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloropropene | | ND | ND | ug/l | NC | 20 |
| 2,2-Dichloropropane | | ND | ND | ug/l | NC | 20 |
| 1,1,1,2-Tetrachloroethane | | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichloropropane | | ND | ND | ug/l | NC | 20 |
| | | | | | | |



Lab Duplicate Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

| Volatile Organics by GC/MS 524.2 Associated sample(s): 01,04-08 QC Batch ID: WG319665-2 QC Sample: L0805753-02 Client ID: D Bromochloromethane ND ND ug/l NC n-Butylbenzene ND ND ug/l NC Dichlorodifluoromethane ND ND ug/l NC Hexachlorobutadiene ND ND ND ug/l NC Isopropylbenzene ND ND ND ug/l NC P-Isopropyltoluene ND ND ND ug/l NC Naphthalene ND ND ND ug/l NC n-Propylbenzene ND ND ND ug/l NC sec-Butylbenzene ND ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ND ug/l NC 1,2,4-Trimethylbenzene ND ND ND ug/l NC | 20 20 20 20 20 20 20 20 |
|--|--|
| n-Butylbenzene ND ND ug/l NC Dichlorodifluoromethane ND ND ug/l NC Hexachlorobutadiene ND ND ug/l NC Isopropylbenzene ND ND ug/l NC p-Isopropyltoluene ND ND ug/l NC Naphthalene ND ND ug/l NC n-Propylbenzene ND ND ug/l NC sec-Butylbenzene ND ND ug/l NC tert-Butylbenzene ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ND ug/l NC | 20 20 20 20 20 20 20 |
| Dichlorodifluoromethane ND ND ND Ug/l NC Isopropylbenzene ND ND ND ND Ug/l NC Isopropylbenzene ND ND ND Ug/l NC NC Naphthalene ND ND ND Ug/l NC NC Naphthalene ND ND ND Ug/l NC NC NC Naphthalene ND ND ND Ug/l NC NC Sec-Butylbenzene ND ND ND Ug/l NC Sec-Butylbenzene ND ND ND Ug/l NC Itert-Butylbenzene ND ND ND Ug/l NC | 20 20 20 20 20 20 |
| Hexachlorobutadiene ND ND ug/l NC Isopropylbenzene ND ND ug/l NC p-Isopropyltoluene ND ND ug/l NC Naphthalene ND ND ug/l NC n-Propylbenzene ND ND ug/l NC sec-Butylbenzene ND ND ug/l NC tert-Butylbenzene ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ug/l NC | 20 20 20 20 |
| Isopropylbenzene ND ND ug/l NC p-Isopropyltoluene ND ND ND ug/l NC Naphthalene ND ND ND ug/l NC n-Propylbenzene ND ND ND ug/l NC sec-Butylbenzene ND ND ND ug/l NC tert-Butylbenzene ND ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ND ug/l NC | 20 20 20 |
| p-Isopropyltoluene ND ND ug/l NC Naphthalene ND ND ND ug/l NC n-Propylbenzene ND ND ND ug/l NC sec-Butylbenzene ND ND ND ug/l NC tert-Butylbenzene ND ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ND ug/l NC | 20 20 |
| Naphthalene ND ND ND ug/l NC n-Propylbenzene ND ND ND ug/l NC sec-Butylbenzene ND ND ND ug/l NC tert-Butylbenzene ND ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ND ug/l NC NC ND ND ug/l NC NC ND ND ug/l NC NC ND ND ug/l NC | 20 |
| n-Propylbenzene ND ND ug/l NC sec-Butylbenzene ND ND ug/l NC tert-Butylbenzene ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND Ug/l NC | |
| sec-Butylbenzene ND ND ug/l NC tert-Butylbenzene ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ug/l NC | 00 |
| tert-Butylbenzene ND ND ug/l NC 1,2,3-Trichlorobenzene ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ug/l NC | 20 |
| 1,2,3-Trichlorobenzene ND ND ug/l NC 1,2,4-Trichlorobenzene ND ND ug/l NC | 20 |
| 1,2,4-Trichlorobenzene ND ND ug/l NC | 20 |
| | 20 |
| 1,2,4-Trimethylbenzene ND ND ug/l NC | 20 |
| · | 20 |
| 1,3,5-Trimethylbenzene ND ND ug/l NC | 20 |
| Bromobenzene ND ND ug/l NC | 20 |
| o-Chlorotoluene ND ND ug/l NC | 20 |
| p-Chlorotoluene ND ND ug/l NC | 20 |
| Dibromomethane ND ND ug/l NC | 20 |
| 1,2-Dibromoethane ND ND ug/l NC | 20 |



WALPOLE PARK SOUTH Bar

Report Date:

Lab Number: L0806023 **Report Date:** 05/05/08

Parameter Native Sample Duplicate Sample Units **RPD RPD Limits** Volatile Organics by GC/MS 524.2 Associated sample(s): 01,04-08 QC Batch ID: WG319665-2 QC Sample: L0805753-02 Client ID: DUP Sample 1,2-Dibromo-3-chloropropane ND ND NC 20 ug/l 1,3-Dichloropropane 20 ND ND ug/l NC Methyl tert butyl ether ND ND ug/l NC 20

| | | | | | Acceptance | |
|------------------------|-----------|-----------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 115 | | 115 | | 80-120 | |
| 4-Bromofluorobenzene | 87 | | 88 | | 80-120 | |



Project Name:

Project Number:

12700058-003

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

Report Date: 05/05/08

| rameter | Na | ative Sample | Duplicate Sa | mple Units | RPD | RPD Limits |
|--------------------------------|--------------------------|-------------------|--------------|----------------------|----------------|-------------|
| latile Organics by GC/MS 524.2 | Associated sample(s): 02 | 2-03 QC Batch ID: | WG320296-2 | QC Sample: L0806023- | -03 Client ID: | MW-3-042808 |
| Methylene chloride | | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethane | | ND | ND | ug/l | NC | 20 |
| Chloroform | | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride | | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane | | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane | | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane | | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene | | ND | ND | ug/l | NC | 20 |
| Chlorobenzene | | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane | | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane | | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane | | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane | | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene | | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene | | ND | ND | ug/l | NC | 20 |
| Bromoform | | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane | | ND | ND | ug/l | NC | 20 |
| Benzene | | ND | ND | ug/l | NC | 20 |
| Toluene | | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date: 05/05/08

| arameter | Native Sample | Duplicate Sa | ample Units | RPD | RPD Limits |
|---|----------------------|--------------|------------------------|--------------|-------------|
| platile Organics by GC/MS 524.2 Associated sample(s |): 02-03 QC Batch ID | WG320296-2 | QC Sample: L0806023-03 | 3 Client ID: | MW-3-042808 |
| Ethylbenzene | ND | ND | ug/l | NC | 20 |
| p/m-Xylene | ND | ND | ug/l | NC | 20 |
| Chloromethane | ND | ND | ug/l | NC | 20 |
| Bromomethane | ND | ND | ug/l | NC | 20 |
| Vinyl chloride | ND | ND | ug/l | NC | 20 |
| Chloroethane | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethene | ND | ND | ug/l | NC | 20 |
| trans-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| cis-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| Trichloroethene | ND | ND | ug/l | NC | 20 |
| 1,2-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,3-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,4-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| Styrene | ND | ND | ug/l | NC | 20 |
| o-Xylene | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloropropene | ND | ND | ug/l | NC | 20 |
| 2,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| 1,1,1,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichloropropane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L0806023

Report Date: 05/05/08

| arameter | rameter | | Sample | Duplicate Sa | ample Uı | nits | RPD | RPD Limits |
|---------------------------------|-----------------------|-------|--------------|--------------|------------|-------------|------------|-------------|
| olatile Organics by GC/MS 524.2 | Associated sample(s): | 02-03 | QC Batch ID: | WG320296-2 | QC Sample: | L0806023-03 | Client ID: | MW-3-042808 |
| Bromochloromethane | | N | ID | ND | ι | ıg/l | NC | 20 |
| n-Butylbenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| Dichlorodifluoromethane | | N | ID | ND | ι | ıg/l | NC | 20 |
| Hexachlorobutadiene | | N | ID | ND | ι | ıg/l | NC | 20 |
| Isopropylbenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| p-Isopropyltoluene | | N | ID | ND | ι | ıg/l | NC | 20 |
| Naphthalene | | N | ID | ND | ι | ıg/l | NC | 20 |
| n-Propylbenzene | | N | ID | ND | U | ıg/l | NC | 20 |
| sec-Butylbenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| tert-Butylbenzene | | N | ID | ND | U | ıg/l | NC | 20 |
| 1,2,3-Trichlorobenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| 1,2,4-Trichlorobenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| 1,2,4-Trimethylbenzene | | N | ID | ND | ι | ıg/l | NC | 20 |
| 1,3,5-Trimethylbenzene | | N | ID | ND | l | ıg/l | NC | 20 |
| Bromobenzene | | N | ID | ND | l | ıg/l | NC | 20 |
| o-Chlorotoluene | | ND | | ND | l | ıg/l | NC | 20 |
| p-Chlorotoluene | | ND | | ND | l | ıg/l | NC | 20 |
| Dibromomethane | | ND | | ND | ι | ıg/l | NC | 20 |
| 1,2-Dibromoethane | | N | ID | ND | ι | ıg/l | NC | 20 |
| | | | | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L0806023

Report Date:

05/05/08

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|-------------------------------------|------------------|-----------------|----------------|-------------|
| Volatile Organics by GC/MS 524.2 Assoc | iated sample(s): 02-03 QC Batch ID: | WG320296-2 QC S | ample: L0806023 | -03 Client ID: | MW-3-042808 |
| 1,2-Dibromo-3-chloropropane | ND | ND | ug/l | NC | 20 |
| 1,3-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Methyl tert butyl ether | ND | ND | ug/l | NC | 20 |

| | | | | Acceptance | | | |
|------------------------|-----------|-----------|-----------|------------|----------|--|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | | |
| 1,2-Dichlorobenzene-d4 | 112 | | 111 | | 80-120 | | |
| 4-Bromofluorobenzene | 90 | | 88 | | 80-120 | | |



METALS



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 09:10

Client ID: RIZ-10-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ries | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:2 | 8 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.062 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 13:4 | 8 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:2 | 8 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 17:2 | 8 EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 10:12

Client ID: RIZ-9-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:3 | 4 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.021 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 13:5 | 5 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:34 | 4 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 17:59 | 9 EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-03 Date Collected: 04/28/08 11:20

Client ID: MW-3-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:4 | 0 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.010 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 13:5 | 7 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:4 | 0 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 3 EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: L0806023-04 Date Collected: 04/28/08 12:15

Client ID: RIZ-8-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:4 | 5 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.025 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 13:5 | 8 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 22:4 | 5 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 18:0 | 6 EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 12:30

Client ID: RIZ-8S-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:0 | 7 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.054 | | mg/l | 0.010 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 (| 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 14:0 | 0 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:0 | 7 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 0 EPA 3005A | 60,6010B | AI |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 14:36

Client ID: GHC-6-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:1: | 3 EPA 3005A | 64,6020A | BM |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.059 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 14:02 | 2 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:13 | 3 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 18:14 | 4 EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

SAMPLE RESULTS

Lab ID: Date Collected: 04/28/08 14:52

Client ID: MW-2-042808 Date Received: 04/29/08
Sample Location: WALPOLE, MA Field Prep: Field Filtered

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|----------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Dissolved Metals by | y MCP 60 | 00/7000 ser | ies | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:18 | 8 EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.031 | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 14:04 | 4 EPA 7470A | 64,7470A | RC |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 04/30/08 12:00 | 05/01/08 23:18 | 8 EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 18:1 | 7 EPA 3005A | 60,6010B | AI |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|-------------------------|------------------|------------|----------|--------------------|------------------|------------------|----------------------|---------|
| Dissolved Metals by MCP | 6000/7000 series | for sample | e(s): 01 | -07 Batch | : WG320004- | 1 | | |
| Arsenic, Dissolved | ND | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Barium, Dissolved | ND | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | Al |
| Beryllium, Dissolved | ND | mg/l | 0.005 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Cadmium, Dissolved | ND | mg/l | 0.004 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | Al |
| Chromium, Dissolved | ND | mg/l | 0.01 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | Al |
| Lead, Dissolved | ND | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Nickel, Dissolved | ND | mg/l | 0.025 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Selenium, Dissolved | ND | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Silver, Dissolved | ND | mg/l | 0.007 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Vanadium, Dissolved | ND | mg/l | 0.010 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |
| Zinc, Dissolved | ND | mg/l | 0.050 | 1 | 04/30/08 12:00 | 05/02/08 15:41 | 60,6010B | AI |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|-------------------------|--------------------|------------|----------|--------------------|------------------|------------------|----------------------|---------|
| Dissolved Metals by MCF | 9 6000/7000 series | for sample | e(s): 01 | I-07 Batch: | WG320013- | 1 | | |
| Antimony, Dissolved | ND | mg/l | 0.0005 | 1 | 04/30/08 12:00 | 05/01/08 21:28 | 64,6020A | ВМ |
| Thallium, Dissolved | ND | mg/l | 0.0005 | 1 | 04/30/08 12:00 | 05/01/08 21:28 | 64,6020A | ВМ |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | | ilution actor | Date Prepared | Date Analyzed | Analytical Method | |
|------------------------|---------------------|-----------|------------|------------------|------------------|------------------|----------------------|----|
| Dissolved Metals by MC | CP 6000/7000 series | for sampl | e(s): 01-0 | 7 Batch: | : WG320203- | 1 | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 05/01/08 18:00 | 05/02/08 13:25 | 64,7470A | RC |



Project Name:WALPOLE PARK SOUTHLab Number:L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

Method Blank Analysis Batch Quality Control

Prep Information

Digestion Method: EPA 7470A



L0806023

Lab Number:

Lab Control Sample Analysis Batch Quality Control

Project Number: 12700058-003 Report Date: 05/05/08

| Parameter | LCS %Recovery | | LCSD ecovery | %Recovery Limits | RPD | RPD Limits |
|--|-----------------------|-------|-----------------|-----------------------|-----|------------|
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s): | 01-07 | Batch: | WG320004-2 WG320004-3 | | |
| Arsenic, Dissolved | 107 | | 106 | 80-120 | 1 | 20 |
| Barium, Dissolved | 94 | | 92 | 80-120 | 2 | 20 |
| Beryllium, Dissolved | 97 | | 95 | 80-120 | 2 | 20 |
| Cadmium, Dissolved | 108 | | 106 | 80-120 | 2 | 20 |
| Chromium, Dissolved | 95 | | 95 | 80-120 | 0 | 20 |
| Lead, Dissolved | 97 | | 95 | 80-120 | 2 | 20 |
| Nickel, Dissolved | 96 | | 95 | 80-120 | 1 | 20 |
| Selenium, Dissolved | 103 | | 103 | 80-120 | 0 | 20 |
| Silver, Dissolved | 94 | | 92 | 80-120 | 2 | 20 |
| Vanadium, Dissolved | 98 | | 96 | 80-120 | 2 | 20 |
| Zinc, Dissolved | 96 | | 93 | 80-120 | 3 | 20 |
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s): | 01-07 | Batch: | WG320013-2 WG320013-3 | | |
| Antimony, Dissolved | 93 | | 96 | 80-120 | 3 | 20 |
| Thallium, Dissolved | 96 | | 97 | 80-120 | 1 | 20 |
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s): | 01-07 | Batch: | WG320203-2 WG320203-3 | | |
| Mercury, Dissolved | 109 | | 105 | 80-120 | 4 | 20 |



Project Name:

WALPOLE PARK SOUTH

Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal A Absent

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|------------------------------|--------|----|------|------|--------|--|
| L0806023-01A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-01B | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-01C | Plastic 500ml HNO3 preserved | Α | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SE-6010S |
| L0806023-02A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-02B | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-02C | Plastic 500ml HNO3 preserved | Α | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SE-6010S |
| L0806023-03A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-03B | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-03C | Plastic 500ml HNO3 preserved | Α | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SE-6010S |
| L0806023-04A | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-04B | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-04C | Plastic 500ml HNO3 preserved | A | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SE-6010S |
| L0806023-05A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |



Lab Number: L0806023

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 **Report Date:** 05/05/08

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|------------------------------|--------|----|------|------|--------|---|
| L0806023-05B | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-05C | Plastic 500ml HNO3 preserved | A | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SE-6010S |
| L0806023-06A | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-06B | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-06C | Plastic 500ml HNO3 preserved | A | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SB-6010S,MCP-SE-6010S |
| L0806023-07A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-07B | Vial HCl preserved | Α | NA | 4 C | Υ | Absent | 524.2 |
| L0806023-07C | Plastic 500ml HNO3 preserved | A | <2 | 4 C | Y | Absent | MCP-CR-6010S,MCP-AG-6010S,MCP-AS-6010S,MCP-NI-6010S,MCP-PB-6010S,MCP-TL-6020S,MCP-V-6010S,MCP-SB-6020S,MCP-7470S,MCP-BE-6010S,MCP-ZN-6010S,MCP-BA-6010S,MCP-SB-6010S,MCP-SE-6010S |
| L0806023-08A | Vial HCI preserved | Α | NA | 4 C | Υ | Absent | 524.2 |

Container Comments

| L0806023-01A | IR Gun |
|--------------|--------|
| L0806023-01B | IR Gun |
| L0806023-01C | IR Gun |
| L0806023-02A | IR Gun |
| L0806023-02B | IR Gun |
| L0806023-02C | IR Gun |
| L0806023-03A | IR Gun |
| L0806023-03B | IR Gun |
| L0806023-03C | IR Gun |



Project Name: WALPOLE PARK SOUTH Lab Number: L0806023

Project Number: 12700058-003 **Report Date:** 05/05/08

Container Information

Container ID Container Type Cooler pH Temp Pres Seal Analysis

Container Comments

| L0806023-04A | IR Gun |
|--------------|--------|
| L0806023-04B | IR Gun |
| L0806023-04C | IR Gun |
| L0806023-05A | IR Gun |
| L0806023-05B | IR Gun |
| L0806023-05C | IR Gun |
| L0806023-06A | IR Gun |
| L0806023-06B | IR Gun |
| L0806023-06C | IR Gun |
| L0806023-07A | IR Gun |
| L0806023-07B | IR Gun |
| L0806023-07C | IR Gun |
| L0806023-08A | IR Gun |



Project Name:WALPOLE PARK SOUTHLab Number:L0806023Project Number:12700058-003Report Date:05/05/08

GLOSSARY

Acronyms

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD- Laboratory Control Sample Duplicate: Refer to LCS.

 MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NI - Not Ignitable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

ND - Not detected at the reported detection limit for the sample.

RDL - Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

The following data qualifiers have been identified for use under the CT DEP Reasonable Confidence Protocols.

- A Spectra identified as "Aldol Condensation Product".
- B The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- J Estimated value. The analyte was tentatively identified; the quantitation is an estimation. (Tentatively identified compounds only.)

Standard Qualifiers

H - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.

Report Format: Data Usability Report



Project Name:WALPOLE PARK SOUTHLab Number:L0806023Project Number:12700058-003Report Date:05/05/08

REFERENCES

- Methods for the Determination of Organic Compounds in Drinking Water Supplement II. EPA/600/R-92/129, August 1992.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). May 2004.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Woods Hole Labs performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Woods Hole Labs be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



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| Billing Information | Deliverables | | Project Information | MANSFIELD, MA | WESTBORO, MA N |
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Report Date: 26-Nov-08 14:04



| \checkmark | Final Report |
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| | Re-Issued Report |
| | Revised Report |

Laboratory Report

Tetra Tech Rizzo One Grant Street - P.O. Box 9005 Framingham, MA 01701

Attn: Ray Johnson

Project: Walpole Park South-Walpole, MA

Project 12700058

| Laboratory ID | Client Sample ID | <u>Matrix</u> | Date Sampled | Date Received |
|----------------------|-------------------|---------------|---------------------|----------------------|
| SA87371-01 | RIZ-10-GW | Ground Water | 11-Nov-08 14:35 | 12-Nov-08 17:15 |
| SA87371-02 | RIZ-8-GW | Ground Water | 11-Nov-08 13:50 | 12-Nov-08 17:15 |
| SA87371-03 | RIZ-3-GW | Ground Water | 11-Nov-08 15:40 | 12-Nov-08 17:15 |
| SA87371-04 | MW-9-GW | Ground Water | 11-Nov-08 11:50 | 12-Nov-08 17:15 |
| SA87371-05 | MW-2-GW | Ground Water | 11-Nov-08 09:50 | 12-Nov-08 17:15 |
| SA87371-06 | RIZ-9-GW | Ground Water | 11-Nov-08 11:00 | 12-Nov-08 17:15 |
| SA87371-07 | MW-3-GW | Ground Water | 11-Nov-08 08:35 | 12-Nov-08 17:15 |
| SA87371-08 | GHC-6-GW | Ground Water | 11-Nov-08 09:20 | 12-Nov-08 17:15 |
| SA87371-09 | Trip Blank 111108 | Aqueous | 11-Nov-08 00:00 | 12-Nov-08 17:15 |

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87600/E87936 Maine # MA138 New Hampshire # 2538 New Jersey # MA011/MA012 New York # 11393/11840 Pennsylvania # 68-04426/68-02924

Rhode Island # 98 USDA # S-51435

Vermont # VT-11393



Authorized by:

Hanibal C. Tayeh, Ph.D.
President/Laboratory Director

Technical Reviewer's Initial:



Spectrum Analytical holds certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please note that this report contains 43 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supercedes any previously dated reports for the laboratory ID(s) referenced above. Where this report

indicated as revised, this report supercedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report is available upon request. This report may not be reproduced, except in full, without written approval from Spectrum Analytical, Inc.

Spectrum Analytical, Inc. is a NELAC accredited laboratory organization and meets NELAC testing standards. Use of the NELAC logo however does not insure that Spectrum is currently accredited for the specific method or analyte indicated. Please refer to our "Quality" web page at www.spectrum-analytical.com for a full listing of our current certifications and fields of accreditation. States in which Spectrum Analytical, Inc. holds NELAC certification are New York, New Hampshire, New Jersey and Florida. All analytical work for Volatile Organic and Air analysis are transferred to and conducted at our 830 Silver Street location (NY-11840, FL-E87936 and NJ-MA012).

CASE NARRATIVE:

The samples were received 3.9 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 2.0 degrees Celsius was used immediately upon receipt of the samples.

MADEP has published a list of analytical methods (CAM) which provides a series of recommended protocols for the acquisition, analysis and reporting of analytical data in support of MCP decisions. "Presumptive Certainty" can be established only for those methods published by the MADEP in the MCP CAM. The compounds and/or elements reported were specifically requested by the client on the Chain of Custody and in some cases may not include the full analyte list as defined in the method.

According to WSC-CAM 5/2004 Rev.4, Table 11 A-1, recovery for some VOC analytes have been deemed potentially difficult. Although they may still be within the recommended 70%-130% recovery range, a range has been set based on historical control limits.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

EPA 624

Laboratory Control Samples:

8111076-BSD1

The RPD result exceeded the QC control limits; however, both percent recoveries were acceptable. Sample results for the QC batch were accepted based on percent recoveries and completeness of QC data.

1,1,1-Trichloroethane

1,1-Dichloroethene

Carbon tetrachloride

Trichlorofluoromethane (Freon 11)

Vinyl chloride

8111194-BSD1

Analyte out of acceptance range.

2-Hexanone (MBK)

Vinyl chloride

EPA 624

Spikes:

8111194-MS1 Source: SA87191-03

The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.

1,1-Dichloroethene

Benzene

Tetrachloroethene

trans-1,2-Dichloroethene

Trichloroethene

8111194-MSD1 Source: SA87191-03

Analyte out of acceptance range.

Toluene

The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.

1,1-Dichloroethene

Benzene

Tetrachloroethene

trans-1,2-Dichloroethene

Trichloroethene

8111291-MS1 Source: SA87501-01

The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.

Carbon tetrachloride

8111291-MSD1 Source: SA87501-01

Analyte out of acceptance range.

1,1,2,2-Tetrachloroethane

1,1,2-Trichloroethane

1,2-Dichlorobenzene

1,3-Dichlorobenzene

1,4-Dichlorobenzene

Benzene

Bromodichloromethane

Bromoform

Chlorobenzene

Chloroform

cis-1,3-Dichloropropene

Dibromochloromethane

Ethylbenzene

Toluene

trans-1,3-Dichloropropene

The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.

Carbon tetrachloride

SW846 6010B

SW846 6010B

Duplicates:

8111042-DUP1 Source: SA87401-01

Analyses are not controlled on RPD values from sample concentrations that are less than 5 times the reporting level. The batch is accepted based upon the difference between the sample and duplicate is less than or equal to the reporting limit.

Lead

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cei |
|------------|------------------------------------|----------------|------|--------------|------------|----------|-----------------|-------------|-------------|---------|----------|
| olatile O | organic Compounds | | | | | | | | | | |
| olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | by method SW846 5030 Water N | ЛS | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111077 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | u u | " | " | " | Х |
| 3-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | X |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | н | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | H . | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | п | " | u u | " | Х |
| 11-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | и | " | " | " | Х |
| 7-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | п | " | " | " | > |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | | | " | > |
| 6-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | | | " | |
| 6-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| -87-5 | | BRL | | μg/l | 1.0 | 1 | " | " | " | | > |
| | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | п | | | | , > |
| | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | п | | | | > |
| | trans-1,3-Dichloropropene | BRL | | | 1.0 | 1 | " | " | " | | <i>,</i> |
| 0-41-4 | Ethylbenzene | | | μg/l | | • | " | | ,, | | |
| 11-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | " | | | | |
| 34-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | | | |
| 10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l " | 10.0 | 1 | | | | | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | | | | | X |
| 0-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | | | | | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | " | | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 8-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| I-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | п | " | " | " | X |
| 79601-23-1 | m,p-Xylene | BRL | | μg/l | 2.0 | 1 | " | " | " | " | X |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| urrogate | recoveries: | | | | | | | | | | |
| 50-00-4 | 4-Bromofluorobenzene | 105 | | 70-13 | 30 % | | " | " | " | " | |
| | Toluene-d8 | 100 | | 70-13 | | | " | " | " | " | |
| | 1,2-Dichloroethane-d4 | 107 | | 70-13 | | | " | u u | " | " | |
| | Dibromofluoromethane | 106 | | 70-13 | | | п | " | " | | |
| | letals by EPA 200/6000 Series Meth | | | - /- | | | | | | | |
| | Filtration | Field Filtered | i | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111039 | |
| oluble M | letals by EPA 6000/7000 Series Met | | • | 1 | | • | | | | | |
| | | | | ua/l | 5 0 | 1 | SW846 6010B | 21 Nov 00 | 21-Nov-08 | Q111040 | |
| 140-22-4 | Silver | BRL BRL | | μg/l μg/l | 5.0 4.0 | 1 | 300040 00101 | ∠ I-INUV-U8 | ∠ 1-INUV-U8 | " | |
| | Arsenic | | | | | | | | | | |

Sample Identification RIZ-10-GW SA87371-01

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 14:35

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|-------------------------------|------------|------|-------|------|----------|-----------------|-----------|-------------|--------|-------|
| Soluble M | letals by EPA 6000/7000 Serie | es Methods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 8 | 111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | u u | | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | " | " | " | " | |
| 7440-28-0 | Thallium | 11.6 | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 36.3 | | μg/l | 7.5 | 1 | u u | | " | " | |
| Soluble M | letals by EPA 200 Series Metl | hods | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 8 | 111043 | Χ |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 13:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|------------|--|---------------|------|-----------|------|----------|-------------------|-------------|-------------------------|---------|-------|
| Volatile (| Organic Compounds | | | | | | | | | | |
| /olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | by method SW846 5030 Water N | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 18-Nov-08 | 18-Nov-08 | 8111291 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Χ |
| 8-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Χ |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 4-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Χ |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 41-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 07-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 56-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | |
| 56-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 8-87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Х |
| 0061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 00-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 91-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | W . | " | " | " | |
| 634-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 08-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | W . | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | u u | " | u u | u u | Χ |
| 00-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 08-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | u u | u u | Χ |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 79601-23- | ¹ m,p-Xylene | BRL | | μg/l | 2.0 | 1 | W . | " | " | " | Х |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| Surrogate | recoveries: | | | | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 86 | | 70-13 | 30 % | | " | " | " | " | |
| | Toluene-d8 | 97 | | 70-13 | | | " | " | " | " | |
| | 1,2-Dichloroethane-d4 | 107 | | 70-13 | | | " | " | " | " | |
| | Dibromofluoromethane | 112 | | 70-13 | | | " | " | " | " | |
| | Ietals by EPA 200/6000 Series Meth | | | | | | | | | | |
| | • | Field Filtere | d | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111030 | |
| olubla N | Filtration Ietals by EPA 6000/7000 Series Met | | u | 11// | | ' | L. 7. 200.770003A | 1-7 1404-00 | . - 140 V-00 | 0111009 | |
| | • | | | | F 0 | 4 | CM046 C040D | 04 Nav. 00 | 04 Nav. 00 | 0111040 | |
| 440-22-4 | | BRL | | µg/l | 5.0 | 1 | SW846 6010B | ∠1-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l " | 4.0 | 1 | | | - | | |
| 440-39-3 | Barium | 27.6 | | μg/l | 5.0 | 1 | " | " | " | " | |

Sample Identification RIZ-8-GW SA87371-02

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 13:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|------------------------------------|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble N | Metals by EPA 6000/7000 Series Met | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | " | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 26.4 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble M | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Х |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 15:40

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|------------|--|---------------|------|-----------|------|----------|-------------------|-------------|-------------------------|---------|-------|
| Volatile C | Organic Compounds | | | | | | <u> </u> | | | | |
| /olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | by method SW846 5030 Water N | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111077 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 8-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | u u | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | u u | " | " | " | Х |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 4-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 41-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 07-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 56-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | |
| 56-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 8-87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 0061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 00-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 91-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | W . | " | " | " | |
| 634-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 08-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | W . | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | W . | " | " | " | Х |
| 00-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 08-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| 79601-23- | ¹ m,p-Xylene | BRL | | μg/l | 2.0 | 1 | W . | " | " | " | Х |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | W . | " | " | " | Х |
| Surrogate | recoveries: | | | | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 104 | | 70-13 | 30 % | | " | " | " | " | |
| | Toluene-d8 | 99 | | 70-13 | | | " | " | " | " | |
| | 1,2-Dichloroethane-d4 | 114 | | 70-13 | | | " | " | " | " | |
| | Dibromofluoromethane | 109 | | 70-13 | | | " | " | " | | |
| | Ietals by EPA 200/6000 Series Meth | | | | * | | | | | | |
| | • | Field Filtere | Н | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111030 | |
| olubla M | Filtration Ietals by EPA 6000/7000 Series Met | | u | 11/7 | | ' | L. 7. 200.770003A | 1-7 1404-00 | . - 140 4-00 | 5111009 | |
| | • | | | | F 0 | 4 | CM046 C040D | 04 Nav. 00 | 04 Nav. 00 | 0111040 | |
| 440-22-4 | | BRL | | µg/l | 5.0 | 1 | SW846 6010B | ∠1-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l " | 4.0 | 1 | | | - | | |
| 440-39-3 | Barium | 86.2 | | μg/l | 5.0 | 1 | " | " | " | " | |

Sample Identification RIZ-3-GW SA87371-03

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 15:40

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|---|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble N | Metals by EPA 6000/7000 Series Me | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | " | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | 24-Nov-08 | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | 21-Nov-08 | " | |
| 7440-66-6 | Zinc | 21.0 | | μg/l | 7.5 | 1 | " | | " | " | |
| Soluble N | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | X |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 11:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|------------|--|----------------|------|--------|------|----------|------------------|--------------------------|--------------------------|---------|-------|
| Volatile (| Organic Compounds | | | | | | <u> </u> | | | | |
| /olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | by method SW846 5030 Water M | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111077 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | п | · · | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Χ |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | п | · · | " | " | Х |
| 8-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | H . | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | п | · · | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | п | " | " | " | Х |
| 7-66-3 | Chloroform | 2.2 | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 4-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | H . | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 41-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 07-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 56-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | п | " | " | " | |
| 56-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 3-87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 0061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 00-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 91-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 634-04-4 | Methyl tert-butyl ether | 2.7 | | μg/l | 1.0 | 1 | n . | " | " | " | |
| 08-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | u u | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | " | " | " | " | Х |
| 00-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | Х |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 79601-23- | ¹ m,p-Xylene | BRL | | μg/l | 2.0 | 1 | n . | " | " | " | Х |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | u | " | " | " | Х |
| Surrogata | recoveries: | | | * | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 103 | | 70-13 | 30 % | | " | " | " | | |
| | Toluene-d8 | 100 | | 70-13 | | | " | " | " | | |
| | 1,2-Dichloroethane-d4 | 113 | | 70-13 | | | " | u u | " | " | |
| | Dibromofluoromethane | 110 | | 70-13 | | | | " | " | " | |
| | Ietals by EPA 200/6000 Series Meth | | | .0 10 | | | | | | | |
| JIUDIC IV | · | Field Filtered | 4 | N/A | | 1 | EPA 200.7/3005A | 14-Nov 09 | 14-Nov 09 | 8111030 | |
| olubla N | Filtration Letals by EPA 6000/7000 Series Met | | u | 111/73 | | ı | LI A 200.1/3003A | 1 -1 -1104-00 | 1 -1 -1104-00 | 0111039 | |
| | Ietals by EPA 6000/7000 Series Met | | | | | 4 | 014104000405 | 04 N= 05 | 04 N 00 | 0444040 | |
| 440-22-4 | | BRL | | μg/l | 5.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l | 4.0 | 1 | " | " | " | | |
| 440-39-3 | Barium | 18.6 | | μg/l | 5.0 | 1 | " | " | " | " | |

Sample Identification MW-9-GW SA87371-04

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 11:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|-----------------------------------|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble M | 1etals by EPA 6000/7000 Series Me | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | u | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 34.7 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble M | 1 EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Х |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 09:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cer |
|---------------------|------------------------------------|----------------|------|----------|------|----------|-----------------|-----------|-----------|---------|-----|
| Volatile C | Organic Compounds | | | | | | | | | | |
| olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | d by method SW846 5030 Water N | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111077 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | X |
| 3-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 11-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 7-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | X |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | Х |
| 6-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | |
| 6-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | X |
| -87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | n n | " | " | " | > |
| 061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | > |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | X |
| 0-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | n n | " | " | " | X |
| 1-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | II . | " | " | " | |
| 34-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 8-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | n n | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | u u | " | " | " | X |
| 0-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | |
| -34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | n n | " | " | " | X |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 8-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | | " | " | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | | " | " | Х |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | II . | " | " | " | X |
| | ¹m,p-Xylene | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| urronate | recoveries: | | | <u> </u> | | | | | | | |
| urrogate 60-00-4 | 4-Bromofluorobenzene | 104 | | 70-13 | 30 % | | " | " | " | " | |
| | Toluene-d8 | 99 | | 70-13 | | | " | " | " | | |
| | 1,2-Dichloroethane-d4 | 111 | | 70-13 | | | " | | " | | |
| | Dibromofluoromethane | 109 | | 70-13 | | | " | | | " | |
| | Tetals by EPA 200/6000 Series Meth | | | , , , | /• | | | | | | |
| 11. | Filtration | Field Filtered | l | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111039 | |
| oluble M | Tetals by EPA 6000/7000 Series Met | | | | | • | | | 2. 30 | | |
| 40-22-4 | Silver | BRL | | μg/l | 5.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l | 4.0 | 1 | " | " | " | " | |
| | AISCHIU | DIVE | | M3'' | 7.0 | | | | | | |

Sample Identification MW-2-GW SA87371-05

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 09:50

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|------------------------------------|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble N | Metals by EPA 6000/7000 Series Met | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | u | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | u | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | u | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 28.0 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble N | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Х |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 11:00

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cer |
|----------------|------------------------------------|----------------|------|----------------|------|----------|-----------------|-------------|--------------|---------|-----|
| Volatile C | Organic Compounds | | | | | | | | | | |
| olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | l by method SW846 5030 Water I | MS | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111077 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | X |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | п | " | " | " | Х |
| 1-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | п | " | " | " | Х |
| 3-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 8-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| ' -66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 6-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| -34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| 7-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| -35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | W | " | " | " | > |
| 6-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | W | " | " | " | |
| 6-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | н | " | " | " | > |
| -87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | H . | " | " | " | > |
| 061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | H . | " | " | " | > |
| 061-02-6 | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | н | " | " | " | > |
| 0-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | H . | " | " | " | > |
| 1-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | H . | " | " | " | |
| 34-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | W | " | " | " | |
| 8-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | п | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | п | " | " | " | X |
| 0-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | п | " | " | " | |
| -34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | п | " | " | " | > |
| 7-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 8-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | | " | " | Х |
| -55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | | " | " | > |
| -01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | | " | " | > |
| -69-4 | Trichlorofluoromethane (Freon 11 | | | μg/l | 1.0 | 1 | " | " | " | " | > |
| -01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | п | " | " | " | > |
| | ¹ m,p-Xylene | BRL | | μg/l | 2.0 | 1 | п | " | " | " | X |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | " | | " | " | × |
| urrogato | recoveries: | | | | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 101 | | 70-13 | 80 % | | " | " | " | | |
| | Toluene-d8 | 99 | | 70-13 | | | " | " | " | | |
| | 1.2-Dichloroethane-d4 | 111 | | 70-13 70-13 | | | " | u u | " | " | |
| | Dibromofluoromethane | 109 | | 70-13 | | | " | | | " | |
| | Ietals by EPA 200/6000 Series Meth | | | .0 70 | | | | | | | |
| | Filtration | Field Filtered | d | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111039 | |
| oluble N | Tetals by EPA 6000/7000 Series Me | | - | * | | • | | | | | |
| | • | BRL | | μg/l | 5.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 40 22 4 | | | | μq/I | 5.0 | 1 | 344040 00100 | ∠ 1-1NUV-UO | Z I TINUV-UO | 0111042 | |
| 140-22-4 | Silver Arsenic | BRL | | μg/l | 4.0 | 1 | " | | | " | |

Sample Identification RIZ-9-GW SA87371-06

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 11:00

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|---|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble N | 1etals by EPA 6000/7000 Series Me | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | u | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | u | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | 24-Nov-08 | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | 21-Nov-08 | " | |
| 7440-66-6 | Zinc | 20.0 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble N | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Χ |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 08:35

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cer |
|------------|------------------------------------|----------------|------|-----------|------|----------|-----------------|-----------|-----------|---------|-----|
| Volatile C | Organic Compounds | | | | | | | | | | |
| olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | d by method SW846 5030 Water N | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 17-Nov-08 | 17-Nov-08 | 8111194 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | X |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | X |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 3-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | n n | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 7-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 6-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 6-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | n . | " | " | " | > |
| -87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | > |
| 10-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | " | | " | " | > |
| 1-78-6 | • | BRL | | μg/l | 10.0 | 1 | " | | | | , |
| 34-04-4 | 2-Hexanone (MBK) | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 8-10-1 | Methyl tert-butyl ether | BRL | | μg/l | 10.0 | 1 | " | | | | |
| i-09-2 | 4-Methyl-2-pentanone (MIBK) | BRL | | | 10.0 | 1 | | " | " | | > |
| | Methylene chloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 0-42-5 | Styrene | | | μg/l | | 1 | " | " | " | " | , |
|)-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | | , | , | | | X |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | | | | | X |
| 8-88-3 | Toluene | BRL | | μg/l " | 1.0 | 1 | | | | | X |
| -55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | | | | | > |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | | | | | X |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | • | " | " | " | > |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | X |
| 79601-23- | -1m,p-Xylene | BRL | | μg/l | 2.0 | 1 | " | " | " | " | X |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| urrogate | recoveries: | | | | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 81 | | 70-13 | 30 % | | " | " | " | " | |
| 037-26-5 | Toluene-d8 | 95 | | 70-13 | 80 % | | " | " | " | " | |
| 7060-07-0 | 1,2-Dichloroethane-d4 | 127 | | 70-13 | 80 % | | u u | " | " | " | |
| 368-53-7 | Dibromofluoromethane | 119 | | 70-13 | 80 % | | " | " | " | " | |
| oluble N | Metals by EPA 200/6000 Series Meth | ods | | | | | | | | | |
| | Filtration | Field Filtered | | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111039 | |
| oluble N | Metals by EPA 6000/7000 Series Met | hods | | | | | | | | | |
| 40-22-4 | · | BRL | | μg/l | 5.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l | 4.0 | 1 | " | " | " | " | |
| | Barium | 11.4 | | μg/l | 5.0 | 1 | | | | | |

Sample Identification MW-3-GW SA87371-07

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 08:35

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|------------------------------------|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble N | Metals by EPA 6000/7000 Series Met | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | u | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | 8.8 | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | u | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 34.5 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble M | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Х |

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 09:20

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert |
|------------|--|----------------|------|-----------|------|----------|-------------------|-------------|-----------|---------|------|
| Volatile C | Organic Compounds | | | | | | <u> </u> | | | | |
| /olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | by method SW846 5030 Water N | //S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111076 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 8-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 4-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 41-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 07-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 56-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | |
| 56-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 8-87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 0061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Х |
| 00-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 91-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 634-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 08-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | " | " | " | " | Х |
| 00-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | u u | " | " | " | Χ |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 79601-23- | ¹ m,p-Xylene | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| Surrogate | recoveries: | | | | | | | | | | |
| 60-00-4 | 4-Bromofluorobenzene | 84 | | 70-13 | 30 % | | " | | " | | |
| | Toluene-d8 | 95 | | 70-13 | | | " | " | " | | |
| | 1,2-Dichloroethane-d4 | 127 | | 70-13 | | | u u | " | " | " | |
| | Dibromofluoromethane | 120 | | 70-13 | | | u u | " | " | " | |
| | Ietals by EPA 200/6000 Series Meth | | | | | | | | | | |
| | • | Field Filtered | d | N/A | | 1 | EPA 200.7/3005A | 14-Nov-08 | 14-Nov-08 | 8111030 | |
| olubla N | Filtration Ietals by EPA 6000/7000 Series Met | | u | 11// | | ' | L. 7. 200.770003A | 1-7 1404-00 | 141404-00 | 5111009 | |
| | • | | | /! | F 0 | 4 | C/V/046 6040D | 21 Nov 00 | 21 Nov 00 | 0111040 | |
| 440-22-4 | | BRL | | µg/l | 5.0 | 1 | SW846 6010B | ∠1-Nov-08 | 21-Nov-08 | 8111042 | |
| | Arsenic | BRL | | μg/l " | 4.0 | 1 | | | - | | |
| 440-39-3 | Barium | 36.8 | | μg/l | 5.0 | 1 | " | " | " | " | |

Sample Identification GHC-6-GW SA87371-08

Client Project # 12700058

<u>Matrix</u> Ground Water Collection Date/Time 11-Nov-08 09:20

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cert. |
|-----------|---|--------|------|-------|------|----------|-----------------|-----------|-----------|---------|-------|
| Soluble M | Metals by EPA 6000/7000 Series Me | thods | | | | | | | | | |
| 7440-41-7 | Beryllium | BRL | | μg/l | 2.0 | 1 | SW846 6010B | 21-Nov-08 | 21-Nov-08 | 8111042 | |
| 7440-43-9 | Cadmium | BRL | | μg/l | 2.5 | 1 | " | " | " | " | |
| 7440-47-3 | Chromium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-02-0 | Nickel | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7439-92-1 | Lead | BRL | | μg/l | 7.5 | 1 | " | " | " | " | |
| 7440-36-0 | Antimony | BRL | | μg/l | 6.0 | 1 | " | " | " | " | |
| 7782-49-2 | Selenium | BRL | | μg/l | 15.0 | 1 | " | " | " | " | |
| 7440-28-0 | Thallium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-62-2 | Vanadium | BRL | | μg/l | 5.0 | 1 | " | " | " | " | |
| 7440-66-6 | Zinc | 21.6 | | μg/l | 7.5 | 1 | " | " | " | " | |
| Soluble M | Metals by EPA 200 Series Methods | | | | | | | | | | |
| 7439-97-6 | Mercury | BRL | | μg/l | 0.20 | 1 | EPA 245.1/7470A | 21-Nov-08 | 24-Nov-08 | 8111043 | Х |

Client Project # 12700058

Matrix Aqueous Collection Date/Time 11-Nov-08 00:00

| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | Dilution | Method Ref. | Prepared | Analyzed | Batch | Cer |
|----------------------|-----------------------------------|-----------|------|-------|-------|----------|-------------|-----------|-----------|---------|-----|
| olatile C | Organic Compounds | | | | | | | | | | |
| /olatile C | Organic Compounds by GCMS | | | | | | | | | | |
| repared | l by method SW846 5030 Water M | 1S | | | | | | | | | |
| 7-64-1 | Acetone | BRL | | μg/l | 20.0 | 1 | EPA 624 | 14-Nov-08 | 15-Nov-08 | 8111076 | |
| 1-43-2 | Benzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-27-4 | Bromodichloromethane | BRL | | μg/l | 1.0 | 1 | " | u u | " | " | Х |
| 5-25-2 | Bromoform | BRL | | μg/l | 1.0 | 1 | " | u u | " | " | Х |
| 4-83-9 | Bromomethane | BRL | | μg/l | 2.0 | 1 | " | u u | " | " | Х |
| 8-93-3 | 2-Butanone (MEK) | BRL | | μg/l | 10.0 | 1 | " | u u | " | " | |
| 6-23-5 | Carbon tetrachloride | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-90-7 | Chlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-00-3 | Chloroethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 7-66-3 | Chloroform | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 4-87-3 | Chloromethane | BRL | | μg/l | 2.0 | 1 | " | " | " | " | Х |
| 24-48-1 | Dibromochloromethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-50-1 | 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 41-73-1 | 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 06-46-7 | 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-34-3 | 1,1-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 7-06-2 | 1,2-Dichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 5-35-4 | 1,1-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 56-59-2 | cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 56-60-5 | trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 3-87-5 | 1,2-Dichloropropane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 0061-01-5 | cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| | trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 00-41-4 | Ethylbenzene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 91-78-6 | 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 634-04-4 | Methyl tert-butyl ether | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 08-10-1 | 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | 1 | " | " | " | " | |
| 5-09-2 | Methylene chloride | BRL | | μg/l | 10.0 | 1 | " | " | " | " | Х |
| 00-42-5 | Styrene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | |
| 9-34-5 | 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 27-18-4 | Tetrachloroethene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 08-88-3 | Toluene | BRL | | μg/l | 1.0 | 1 | " | " | " | " | Х |
| 1-55-6 | 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | 1 | " | " | " | | Х |
| 9-00-5 | 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | 1 | | " | " | | X |
| 9-01-6 | Trichloroethene | BRL | | μg/l | 1.0 | 1 | | " | " | | Х |
| 5-69-4 | Trichlorofluoromethane (Freon 11) | | | μg/l | 1.0 | 1 | | " | " | " | Х |
| 5-01-4 | Vinyl chloride | BRL | | μg/l | 1.0 | 1 | | " | " | " | X |
| | 1m,p-Xylene | BRL | | μg/l | 2.0 | 1 | | " | " | " | X |
| 5-47-6 | o-Xylene | BRL | | μg/l | 1.0 | 1 | | " | " | " | X |
| | recoveries: | | | | | | | | | | |
| ourrogate 60-00-4 | 4-Bromofluorobenzene | 82 | | 70-13 | 30 % | | " | " | " | | |
| | Toluene-d8 | 93 | | 70-13 | | | " | | " | | |
| | 1,2-Dichloroethane-d4 | 93 127 | | 70-13 | | | " | | " | | |
| | | 161 | | 10-1 | JU /0 | | | | | | |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|--|--------------|------|--------------|------|----------------|------------------|-----------|------------------|-----|--------------|
| Batch 8111076 - SW846 5030 Water MS | | | | | | | | | | |
| Blank (8111076-BLK1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Acetone | BRL | | μg/l | 20.0 | | | | | | |
| Benzene | BRL | | μg/l | 1.0 | | | | | | |
| Bromodichloromethane | BRL | | μg/l | 1.0 | | | | | | |
| Bromoform | BRL | | μg/l | 1.0 | | | | | | |
| Bromomethane | BRL | | μg/l | 2.0 | | | | | | |
| 2-Butanone (MEK) | BRL | | μg/l | 10.0 | | | | | | |
| Carbon tetrachloride | BRL | | μg/l | 1.0 | | | | | | |
| Chlorobenzene | BRL | | μg/l | 1.0 | | | | | | |
| Chloroethane | BRL | | μg/l | 2.0 | | | | | | |
| Chloroform | BRL | | μg/l | 1.0 | | | | | | |
| Chloromethane | BRL | | μg/l | 2.0 | | | | | | |
| Dibromochloromethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,2-Dichlorobenzene | BRL | | μg/l | 1.0 | | | | | | |
| 1,3-Dichlorobenzene | BRL | | μg/l | 1.0 | | | | | | |
| 1,4-Dichlorobenzene | BRL | | μg/l | 1.0 | | | | | | |
| 1,1-Dichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,2-Dichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,1-Dichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| cis-1,2-Dichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| trans-1,2-Dichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| 1,2-Dichloropropane | BRL | | μg/l | 1.0 | | | | | | |
| cis-1,3-Dichloropropene | BRL | | μg/l | 1.0 | | | | | | |
| trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | | | | | | |
| Ethylbenzene | BRL | | μg/l | 1.0 | | | | | | |
| 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | | | | | | |
| Methyl tert-butyl ether | BRL | | μg/l | 1.0 | | | | | | |
| 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | | | | | | |
| Methylene chloride | BRL | | μg/l | 10.0 | | | | | | |
| Styrene | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | | | | | | |
| Tetrachloroethene | BRL | | μg/l | 1.0 | | | | | | |
| Toluene | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| Trichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | | | | | | |
| Vinyl chloride | BRL | | μg/l | 1.0 | | | | | | |
| m,p-Xylene | BRL | | μg/l | 2.0 | | | | | | |
| o-Xylene | BRL | | μg/l | 1.0 | | | | | | |
| Surrogate: 4-Bromofluorobenzene | 25.2 | | μg/l | | 30.0 | | 84 | 70-130 | | |
| Surrogate: Toluene-d8 Surrogate: 1,2-Dichloroethane-d4 | 29.2 33.0 | | μg/l μg/l | | 30.0 30.0 | | 98 110 | 70-130 70-130 | | |
| Surrogate: Dibromofluoromethane | 32.3 | | μg/l | | 30.0 | | 108 | 70-130 70-130 | | |
| LCS (8111076-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Acetone | 18.0 | | μg/l | | 20.0 | | 90 | 70-130 | | |
| Benzene | 21.7 | | μg/l | | 20.0 | | 108 | 70-130 | | |
| Bromodichloromethane | 24.1 | | μg/l | | 20.0 | | 121 | 35-155 | | |
| Bromoform | 17.5 | | μg/l | | 20.0 | | 87 | 45-169 | | |
| Bromomethane | 22.2 | | μg/l | | 20.0 | | 111 | 1-242 | | |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|---|--------------|------|--------------|------|----------------|------------------|------------|------------------|-----|--------------|
| Batch 8111076 - SW846 5030 Water MS | | | | | | | | | | |
| LCS (8111076-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| 2-Butanone (MEK) | 18.1 | | μg/l | | 20.0 | | 90 | 70-130 | | |
| Carbon tetrachloride | 23.3 | | μg/l | | 20.0 | | 117 | 70-140 | | |
| Chlorobenzene | 21.2 | | μg/l | | 20.0 | | 106 | 70-130 | | |
| Chloroethane | 22.4 | | μg/l | | 20.0 | | 112 | 14-230 | | |
| Chloroform | 23.0 | | μg/l | | 20.0 | | 115 | 51-138 | | |
| Chloromethane | 20.3 | | μg/l | | 20.0 | | 101 | 1-273 | | |
| Dibromochloromethane | 19.2 | | μg/l | | 20.0 | | 96 | 53-149 | | |
| 1,2-Dichlorobenzene | 20.4 | | μg/l | | 20.0 | | 102 | 18-190 | | |
| 1,3-Dichlorobenzene | 22.2 | | μg/l | | 20.0 | | 111 | 59-156 | | |
| 1,4-Dichlorobenzene | 20.4 | | μg/l | | 20.0 | | 102 | 18-190 | | |
| 1,1-Dichloroethane | 20.8 | | μg/l | | 20.0 | | 104 | 59-155 | | |
| 1,2-Dichloroethane | 20.1 | | μg/l | | 20.0 | | 100 | 49-155 | | |
| 1,1-Dichloroethene | 20.1 | | μg/l | | 20.0 | | 100 | 70-130 | | |
| cis-1,2-Dichloroethene | 20.0 | | μg/l | | 20.0 | | 100 | 70-130 | | |
| trans-1,2-Dichloroethene | 21.0 | | μg/l | | 20.0 | | 105 | 54-156 | | |
| 1,2-Dichloropropane | 21.4 | | μg/l | | 20.0 | | 107 | 1-210 | | |
| cis-1,3-Dichloropropene | 18.6 | | μg/l | | 20.0 | | 93 | 1-227 | | |
| trans-1,3-Dichloropropene | 18.3 | | μg/l | | 20.0 | | 92 | 17-183 | | |
| Ethylbenzene | 22.0 | | μg/l | | 20.0 | | 110 | 37-162 | | |
| 2-Hexanone (MBK) | 14.5 | | μg/l | | 20.0 | | 72 | 70-130 | | |
| Methyl tert-butyl ether | 20.2 | | μg/l | | 20.0 | | 101 | 70-130 | | |
| 4-Methyl-2-pentanone (MIBK) | 18.1 | | μg/l | | 20.0 | | 90 | 70-130 | | |
| Methylene chloride | 22.5 | | μg/l | | 20.0 | | 113 | 1-221 | | |
| Styrene | 19.4 | | μg/l | | 20.0 | | 97 | 70-130 | | |
| 1,1,2,2-Tetrachloroethane | 20.1 | | | | 20.0 | | 101 | 46-157 | | |
| Tetrachloroethene | 19.0 | | μg/l | | 20.0 | | 95 | 64-148 | | |
| Toluene | 21.0 | | μg/l | | 20.0 | | 105 | 70-130 | | |
| 1,1,1-Trichloroethane | 20.0 | | μg/l | | 20.0 | | 100 | 52-162 | | |
| 1,1,2-Trichloroethane | 20.7 | | µg/l | | 20.0 | | 103 | 52-162 | | |
| Trichloroethene | 20.7 | | μg/l μg/l | | 20.0 | | 105 | 71-157 | | |
| | | | | | | | | | | |
| Trichlorofluoromethane (Freon 11) Vinyl chloride | 21.4 23.8 | | μg/l | | 20.0 20.0 | | 107 119 | 17-181 1-251 | | |
| • | | | µg/l | | | | | | | |
| m,p-Xylene | 45.3 | | μg/l | | 40.0 | | 113 | 70-130 70-130 | | |
| o-Xylene Surrogate: 4-Bromofluorobenzene | 22.8 31.3 | | µg/l | | 20.0 30.0 | | 114 | 70-130 70-130 | | |
| Surrogate: 4-Bromonuoropenzene Surrogate: Toluene-d8 | 31.3 30.2 | | μg/l μg/l | | 30.0 30.0 | | 104 101 | 70-130 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 29.2 | | μg/l | | 30.0 | | 97 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 30.8 | | μg/l | | 30.0 | | 103 | 70-130 | | |
| LCS Dup (8111076-BSD1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Acetone | 17.0 | | μg/l | | 20.0 | | 85 | 70-130 | 6 | 30 |
| Benzene | 17.3 | | μg/l | | 20.0 | | 87 | 70-130 | 22 | 30 |
| Bromodichloromethane | 20.3 | | μg/l | | 20.0 | | 101 | 35-155 | 17 | 30 |
| Bromoform | 16.3 | | μg/l | | 20.0 | | 82 | 45-169 | 7 | 30 |
| Bromomethane | 17.8 | | μg/l | | 20.0 | | 89 | 1-242 | 22 | 30 |
| 2-Butanone (MEK) | 19.1 | | μg/l | | 20.0 | | 96 | 70-130 | 6 | 30 |
| Carbon tetrachloride | 15.8 | QR2 | μg/l | | 20.0 | | 79 | 70-140 | 39 | 30 |
| Chlorobenzene | 17.6 | | μg/l | | 20.0 | | 88 | 70-130 | 18 | 30 |
| Chloroethane | 17.8 | | μg/l | | 20.0 | | 89 | 14-230 | 23 | 30 |
| Chloroform | 19.4 | | μg/l | | 20.0 | | 97 | 51-138 | 17 | 30 |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|-----------------------------------|----------------|------|-------|------|----------------|------------------|------|----------------|-----|--------------|
| Batch 8111076 - SW846 5030 Water | MS | | | | | | | | | |
| LCS Dup (8111076-BSD1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Chloromethane | 15.6 | | μg/l | | 20.0 | | 78 | 1-273 | 26 | 30 |
| Dibromochloromethane | 17.3 | | μg/l | | 20.0 | | 87 | 53-149 | 10 | 30 |
| 1,2-Dichlorobenzene | 17.4 | | μg/l | | 20.0 | | 87 | 18-190 | 16 | 30 |
| 1,3-Dichlorobenzene | 19.3 | | μg/l | | 20.0 | | 96 | 59-156 | 14 | 30 |
| 1,4-Dichlorobenzene | 17.5 | | μg/l | | 20.0 | | 88 | 18-190 | 15 | 30 |
| 1,1-Dichloroethane | 17.3 | | μg/l | | 20.0 | | 87 | 59-155 | 18 | 30 |
| 1,2-Dichloroethane | 17.9 | | μg/l | | 20.0 | | 90 | 49-155 | 11 | 30 |
| 1,1-Dichloroethene | 14.2 | QR2 | μg/l | | 20.0 | | 71 | 70-130 | 34 | 30 |
| cis-1,2-Dichloroethene | 16.8 | | μg/l | | 20.0 | | 84 | 70-130 | 18 | 30 |
| trans-1,2-Dichloroethene | 16.4 | | μg/l | | 20.0 | | 82 | 54-156 | 24 | 30 |
| 1,2-Dichloropropane | 17.9 | | μg/l | | 20.0 | | 89 | 1-210 | 18 | 30 |
| cis-1,3-Dichloropropene | 16.1 | | μg/l | | 20.0 | | 80 | 1-227 | 14 | 30 |
| trans-1,3-Dichloropropene | 16.3 | | μg/l | | 20.0 | | 82 | 17-183 | 11 | 30 |
| Ethylbenzene | 17.1 | | μg/l | | 20.0 | | 86 | 37-162 | 25 | 30 |
| 2-Hexanone (MBK) | 14.2 | | μg/l | | 20.0 | | 71 | 70-130 | 2 | 30 |
| Methyl tert-butyl ether | 19.1 | | μg/l | | 20.0 | | 96 | 70-130 | 6 | 30 |
| 4-Methyl-2-pentanone (MIBK) | 18.6 | | μg/l | | 20.0 | | 93 | 70-130 | 3 | 30 |
| Methylene chloride | 19.0 | | μg/l | | 20.0 | | 95 | 1-221 | 17 | 30 |
| Styrene | 15.7 | | μg/l | | 20.0 | | 78 | 70-130 | 21 | 30 |
| 1,1,2,2-Tetrachloroethane | 19.4 | | μg/l | | 20.0 | | 97 | 46-157 | 4 | 30 |
| Tetrachloroethene | 14.4 | | μg/l | | 20.0 | | 72 | 64-148 | 27 | 30 |
| Toluene | 17.0 | | μg/l | | 20.0 | | 85 | 70-130 | 21 | 30 |
| 1,1,1-Trichloroethane | 14.6 | QR2 | μg/l | | 20.0 | | 73 | 52-162 | 31 | 30 |
| 1,1,2-Trichloroethane | 18.8 | | μg/l | | 20.0 | | 94 | 52-150 | 9 | 30 |
| Trichloroethene | 16.3 | | μg/l | | 20.0 | | 82 | 71-157 | 25 | 30 |
| Trichlorofluoromethane (Freon 11) | 13.9 | QR2 | μg/l | | 20.0 | | 69 | 17-181 | 43 | 30 |
| Vinyl chloride | 15.8 | QR2 | μg/l | | 20.0 | | 79 | 1-251 | 41 | 30 |
| m,p-Xylene | 35.6 | | μg/l | | 40.0 | | 89 | 70-130 | 24 | 30 |
| o-Xylene | 18.4 | | μg/l | | 20.0 | | 92 | 70-130 | 21 | 30 |
| Surrogate: 4-Bromofluorobenzene | 30.9 | | μg/l | | 30.0 | | 103 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.4 | | μg/l | | 30.0 | | 101 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 29.8 | | μg/l | | 30.0 | | 99 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 31.1 | | μg/l | | 30.0 | | 104 | 70-130 | | |
| Matrix Spike (8111076-MS1) | Source: SA8726 | 7-01 | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Benzene | 17.5 | | μg/l | | 20.0 | BRL | 88 | 70-130 | | |
| Bromodichloromethane | 23.0 | | μg/l | | 20.0 | BRL | 115 | 35-155 | | |
| Bromoform | 17.2 | | μg/l | | 20.0 | BRL | 86 | 45-169 | | |
| Bromomethane | 11.9 | | μg/l | | 20.0 | BRL | 60 | 1-242 | | |
| Carbon tetrachloride | 21.0 | | μg/l | | 20.0 | BRL | 105 | 70-140 | | |
| Chlorobenzene | 19.3 | | μg/l | | 20.0 | BRL | 97 | 70-130 | | |
| Chloroethane | 13.9 | | μg/l | | 20.0 | BRL | 70 | 14-230 | | |
| Chloroform | 21.5 | | μg/l | | 20.0 | BRL | 108 | 51-138 | | |
| Chloromethane | 10.0 | | μg/l | | 20.0 | BRL | 50 | 1-273 | | |
| Dibromochloromethane | 18.5 | | μg/l | | 20.0 | BRL | 92 | 53-149 | | |
| 1,2-Dichlorobenzene | 21.4 | | μg/l | | 20.0 | BRL | 107 | 18-190 | | |
| 1,3-Dichlorobenzene | 21.8 | | μg/l | | 20.0 | BRL | 109 | 59-156 | | |
| 1,4-Dichlorobenzene | 20.1 | | μg/l | | 20.0 | BRL | 100 | 18-190 | | |
| 1,1-Dichloroethane | 19.3 | | μg/l | | 20.0 | BRL | 97 | 59-155 | | |
| 1,2-Dichloroethane | 18.0 | | μg/l | | 20.0 | BRL | 90 | 49-155 | | |

| Analyta(s) | Popult Floo | Unita | *RDL | Spike | Source Result | %REC | %REC | RPD | RPD Limit |
|--|--------------------|--------------|------|--------------|------------------|-----------|------------------|----------|--------------|
| Analyte(s) | Result Flag | Units | *KDL | Level | Resuit | %KEC | Limits | KPD | Limit |
| Batch 8111076 - SW846 5030 Water | r MS | | | | | | | | |
| Matrix Spike (8111076-MS1) | Source: SA87267-01 | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | |
| 1,1-Dichloroethene | 14.1 | μg/l | | 20.0 | BRL | 71 | 70-130 | | |
| trans-1,2-Dichloroethene | 14.1 | μg/l | | 20.0 | BRL | 71 | 54-156 | | |
| 1,2-Dichloropropane | 20.1 | μg/l | | 20.0 | BRL | 101 | 1-210 | | |
| cis-1,3-Dichloropropene | 16.9 | μg/l | | 20.0 | BRL | 84 | 1-227 | | |
| trans-1,3-Dichloropropene | 17.0 | μg/l | | 20.0 | BRL | 85 | 17-183 | | |
| Ethylbenzene | 19.3 | μg/l | | 20.0 | BRL | 97 | 37-162 | | |
| Methylene chloride | 17.8 | μg/l | | 20.0 | BRL | 89 | 1-221 | | |
| 1,1,2,2-Tetrachloroethane | 20.1 | μg/l | | 20.0 | BRL | 101 | 46-157 | | |
| Tetrachloroethene | 16.0 | μg/l | | 20.0 | BRL | 80 | 64-148 | | |
| Toluene | 17.6 | μg/l | | 20.0 | BRL | 88 | 70-130 | | |
| 1,1,1-Trichloroethane | 19.1 | μg/l | | 20.0 | BRL | 96 | 52-162 | | |
| 1,1,2-Trichloroethane | 20.8 | μg/l | | 20.0 | BRL | 104 | 52-150 | | |
| Trichloroethene | 17.5 | μg/l | | 20.0 | BRL | 88 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 16.6 | μg/l | | 20.0 | BRL | 83 | 17-181 | | |
| Vinyl chloride | 11.4 | μg/l | | 20.0 | BRL | 57 | 1-251 | | |
| Surrogate: 4-Bromofluorobenzene | 30.6 | μg/l | | 30.0 | | 102 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.1 | μg/l | | 30.0 | | 100 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 Surrogate: Dibromofluoromethane | 29.6 31.5 | μg/l μg/l | | 30.0 30.0 | | 99 105 | 70-130 70-130 | | |
| Matrix Spike Dup (8111076-MSD1) | Source: SA87267-01 | F3 | | 00.0 | | 700 | | | |
| Prepared & Analyzed: 14-Nov-08 | Source. SA67267-01 | | | | | | | | |
| Benzene | 17.5 | ua/l | | 20.0 | BRL | 88 | 70-130 | 0.06 | 30 |
| Bromodichloromethane | 23.5 | μg/l | | 20.0 | BRL | | 35-155 | 2 | 30 |
| Bromoform | 23.5 17.2 | μg/l | | 20.0 | BRL | 118 86 | 45-169 | 0.4 | 30 |
| Bromomethane | 12.0 | μg/l | | 20.0 | BRL | 60 | 1-242 | 0.4 | 30 |
| Carbon tetrachloride | 20.9 | μg/l | | 20.0 | BRL | 104 | 70-140 | 0.9 | 30 |
| Chlorobenzene | 19.1 | μg/l | | 20.0 | BRL | 96 | 70-140 70-130 | 0.7 | 30 |
| Chloroethane | 13.0 | µg/l | | 20.0 | BRL | 65 | 14-230 | 7 | 30 |
| Chloroform | 21.5 | μg/l | | 20.0 | BRL | 107 | 51-138 | 0.2 | 30 |
| Chloromethane | 9.4 | μg/l | | 20.0 | BRL | 47 | 1-273 | 7 | 30 |
| | | µg/l | | | | | | | |
| Dibromochloromethane | 19.0 | µg/l | | 20.0 | BRL | 95 105 | 53-149 | 2 | 30 |
| 1,2-Dichlorobenzene | 21.1 21.8 | μg/l | | 20.0 20.0 | BRL BRL | 105 | 18-190 | 1 | 30 30 |
| 1,3-Dichlorobenzene | | μg/l | | | | 109 | 59-156 | 0.2 | |
| 1,4-Dichlorobenzene | 19.8 | μg/l | | 20.0 20.0 | BRL BRL | 99 | 18-190 | 1 | 30 30 |
| 1,1-Dichloroethane | 19.1 18.2 | µg/l | | 20.0 | BRL | 95 01 | 59-155 | 1 | 30 |
| 1,2-Dichloroethane | | μg/l | | 20.0 | | 91 | 49-155 | 1 | 30 |
| 1,1-Dichloroethene | 14.4 | µg/l | | | BRL BRL | 72 71 | 70-130 | 2 | 30 |
| trans-1,2-Dichloroethene 1,2-Dichloropropane | 14.2 19.9 | μg/l | | 20.0 20.0 | BRL | 71 100 | 54-156 1-210 | 0.2 1 | 30 |
| cis-1,3-Dichloropropene | 17.4 | μg/l | | 20.0 | BRL | 87 | 1-210 | | 30 |
| trans-1,3-Dichloropropene | 17.4 17.5 | μg/l | | 20.0 | BRL | 88 | 17-183 | 3 3 | 30 |
| | | µg/l | | | | | | | |
| Ethylbenzene Methylene chloride | 19.8 17.5 | µg/l | | 20.0 | BRL BRL | 99 87 | 37-162 1 221 | 2 | 30 30 |
| Methylene chloride | 17.5 19.6 | µg/l | | 20.0 | BRL | 87 | 1-221 46 157 | 2 | 30 |
| 1,1,2,2-Tetrachloroethane | | μg/l | | 20.0 | | 98 | 46-157 | 3 | |
| Tetrachloroethene | 16.5 | µg/l | | 20.0 | BRL | 83 | 64-148 | 4 | 30 |
| Toluene | 18.1 | μg/l | | 20.0 | BRL | 91 | 70-130 | 3 | 30 |
| 1,1,1-Trichloroethane | 19.0 | μg/l | | 20.0 | BRL | 95 403 | 52-162 | 0.5 | 30 |
| 1,1,2-Trichloroethane | 20.6 | μg/l | | 20.0 | BRL | 103 | 52-150 | 1 | 30 |
| Trichloroethene | 17.7 | μg/l " | | 20.0 | BRL | 88 | 71-157 | 1 | 30 |
| Trichlorofluoromethane (Freon 11) | 15.7 | μg/l | | 20.0 | BRL | 79 | 17-181 | 5 | 30 |

| | | | | Spike | Source | | %REC | | RPD |
|---|--------------------|--------------|------|--------------|--------|------|------------------|-----|-------|
| Analyte(s) | Result Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111076 - SW846 5030 Water | r MS | | | | | | | | |
| Matrix Spike Dup (8111076-MSD1) | Source: SA87267-01 | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | |
| Vinyl chloride | 11.5 | μg/l | | 20.0 | BRL | 58 | 1-251 | 0.6 | 30 |
| Surrogate: 4-Bromofluorobenzene | 30.9 | μg/l | | 30.0 | | 103 | 70-130 | | |
| Surrogate: Toluene-d8 | 31.0 | μg/l | | 30.0 | | 103 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 30.2 | μg/l | | 30.0 | | 101 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 31.4 | μg/l | | 30.0 | | 105 | 70-130 | | |
| Batch 8111077 - SW846 5030 Water | r MS | | | | | | | | |
| Blank (8111077-BLK1) Prepared & Analyzed: 14-Nov-08 | | | | | | | | | |
| Acetone | BRL | ua/l | 20.0 | | | | | | |
| | | μg/l | | | | | | | |
| Benzene | BRL | μg/l | 1.0 | | | | | | |
| Bromodichloromethane | BRL | μg/l | 1.0 | | | | | | |
| Bromoform | BRL | μg/l | 1.0 | | | | | | |
| Bromomethane | BRL | μg/l | 2.0 | | | | | | |
| 2-Butanone (MEK) | BRL | μg/l | 10.0 | | | | | | |
| Carbon tetrachloride | BRL | μg/l | 1.0 | | | | | | |
| Chlorobenzene | BRL | μg/l | 1.0 | | | | | | |
| Chloroethane | BRL | μg/l | 2.0 | | | | | | |
| Chloroform | BRL | μg/l | 1.0 | | | | | | |
| Chloromethane | BRL | μg/l | 2.0 | | | | | | |
| Dibromochloromethane | BRL | μg/l | 1.0 | | | | | | |
| 1,2-Dichlorobenzene | BRL | μg/l | 1.0 | | | | | | |
| 1,3-Dichlorobenzene | BRL | μg/l | 1.0 | | | | | | |
| 1,4-Dichlorobenzene | BRL | μg/l | 1.0 | | | | | | |
| 1,1-Dichloroethane | BRL | μg/l | 1.0 | | | | | | |
| 1,2-Dichloroethane | BRL | μg/l | 1.0 | | | | | | |
| 1,1-Dichloroethene | BRL | μg/l | 1.0 | | | | | | |
| cis-1,2-Dichloroethene | BRL | μg/l | 1.0 | | | | | | |
| trans-1,2-Dichloroethene | BRL | | 1.0 | | | | | | |
| | | μg/l | | | | | | | |
| 1,2-Dichloropropane | BRL | μg/l | 1.0 | | | | | | |
| cis-1,3-Dichloropropene | BRL | μg/l | 1.0 | | | | | | |
| trans-1,3-Dichloropropene | BRL | μg/l | 1.0 | | | | | | |
| Ethylbenzene | BRL | μg/l | 1.0 | | | | | | |
| 2-Hexanone (MBK) | BRL | μg/l | 10.0 | | | | | | |
| Methyl tert-butyl ether | BRL | μg/l | 1.0 | | | | | | |
| 4-Methyl-2-pentanone (MIBK) | BRL | μg/l | 10.0 | | | | | | |
| Methylene chloride | BRL | μg/l | 10.0 | | | | | | |
| Styrene | BRL | μg/l | 1.0 | | | | | | |
| 1,1,2,2-Tetrachloroethane | BRL | μg/l | 1.0 | | | | | | |
| Tetrachloroethene | BRL | μg/l | 1.0 | | | | | | |
| Toluene | BRL | μg/l | 1.0 | | | | | | |
| 1,1,1-Trichloroethane | BRL | μg/l | 1.0 | | | | | | |
| 1,1,2-Trichloroethane | BRL | μg/l | 1.0 | | | | | | |
| Trichloroethene | BRL | μg/l | 1.0 | | | | | | |
| Trichlorofluoromethane (Freon 11) | BRL | μg/l | 1.0 | | | | | | |
| Vinyl chloride | BRL | μg/l | 1.0 | | | | | | |
| m,p-Xylene | BRL | μg/l | 2.0 | | | | | | |
| o-Xylene | BRL | μg/l | 1.0 | | | | | | |
| Surrogate: 4-Bromofluorobenzene | 30.9 | μg/l | 1.0 | 30.0 | | 103 | 70-130 | | |
| Surrogate: 4-Bromonuorobenzene Surrogate: Toluene-d8 | 30.9 29.7 | μg/l μg/l | | 30.0 30.0 | | 99 | 70-130 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 30.9 | μg/l | | 30.0 | | 103 | 70-130 | | |

| | | | | | Spike | Source | | %REC | | RPD |
|---|--------------|------|--------------|------|--------------|--------|------------|------------------|-----|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111077 - SW846 5030 Water MS | | | | | | | | | | |
| Blank (8111077-BLK1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Surrogate: Dibromofluoromethane | 30.5 | | μg/l | | 30.0 | | 102 | 70-130 | | |
| _CS (8111077-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | | |
| Acetone | 19.6 | | μg/l | | 20.0 | | 98 | 70-130 | | |
| Benzene | 21.7 | | μg/l | | 20.0 | | 108 | 70-130 | | |
| Bromodichloromethane | 19.9 | | μg/l | | 20.0 | | 99 | 35-155 | | |
| Bromoform | 18.7 | | μg/l | | 20.0 | | 94 | 45-169 | | |
| Bromomethane | 10.7 | | μg/l | | 20.0 | | 54 | 1-242 | | |
| 2-Butanone (MEK) | 18.7 | | μg/l | | 20.0 | | 93 | 70-130 | | |
| Carbon tetrachloride | 20.3 | | μg/l | | 20.0 | | 102 | 70-140 | | |
| Chlorobenzene | 21.8 | | μg/l | | 20.0 | | 109 | 70-130 | | |
| Chloroethane | 24.6 | | μg/l | | 20.0 | | 123 | 14-230 | | |
| Chloroform | 22.2 | | μg/l | | 20.0 | | 111 | 51-138 | | |
| Chloromethane | 19.5 | | μg/l | | 20.0 | | 98 | 1-273 | | |
| Dibromochloromethane | 20.1 | | μg/l | | 20.0 | | 101 | 53-149 | | |
| 1,2-Dichlorobenzene | 22.0 | | μg/l | | 20.0 | | 110 | 18-190 | | |
| 1,3-Dichlorobenzene | 21.7 | | μg/l | | 20.0 | | 108 | 59-156 | | |
| 1,4-Dichlorobenzene | 21.3 | | μg/l | | 20.0 | | 106 | 18-190 | | |
| 1,1-Dichloroethane | 21.9 | | μg/l | | 20.0 | | 109 | 59-155 | | |
| 1,2-Dichloroethane | 21.6 | | μg/l | | 20.0 | | 108 | 49-155 | | |
| 1,1-Dichloroethene | 21.1 | | μg/l | | 20.0 | | 105 | 70-130 | | |
| cis-1,2-Dichloroethene | 20.5 | | μg/l | | 20.0 | | 103 | 70-130 | | |
| rans-1,2-Dichloroethene | 22.4 | | μg/l | | 20.0 | | 112 | 54-156 | | |
| 1,2-Dichloropropane | 21.5 | | μg/l | | 20.0 | | 108 | 1-210 | | |
| cis-1,3-Dichloropropene | 20.4 | | μg/l | | 20.0 | | 102 | 1-227 | | |
| rans-1,3-Dichloropropene | 20.0 | | μg/l | | 20.0 | | 100 | 17-183 | | |
| Ethylbenzene | 21.2 | | μg/l | | 20.0 | | 106 | 37-162 | | |
| 2-Hexanone (MBK) | 20.0 | | μg/l | | 20.0 | | 100 | 70-130 | | |
| Methyl tert-butyl ether | 21.9 | | μg/l | | 20.0 | | 110 | 70-130 | | |
| 4-Methyl-2-pentanone (MIBK) | 18.7 | | μg/l | | 20.0 | | 94 | 70-130 | | |
| Methylene chloride | 22.2 | | μg/l | | 20.0 | | 111 | 1-221 | | |
| Styrene | 20.6 | | μg/l | | 20.0 | | 103 | 70-130 | | |
| 1,1,2,2-Tetrachloroethane | 20.9 | | μg/l | | 20.0 | | 104 | 46-157 | | |
| Tetrachloroethene | 21.0 | | μg/l | | 20.0 | | 105 | 64-148 | | |
| Toluene | 21.1 | | μg/l | | 20.0 | | 106 | 70-130 | | |
| 1,1,1-Trichloroethane | 21.3 | | μg/l | | 20.0 | | 107 | 52-162 | | |
| 1,1,2-Trichloroethane | 20.2 | | μg/l | | 20.0 | | 101 | 52-150 | | |
| Trichloroethene | 21.0 | | μg/l | | 20.0 | | 105 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 23.3 | | μg/l | | 20.0 | | 116 | 17-181 | | |
| Vinyl chloride | 21.4 | | μg/l | | 20.0 | | 107 | 1-251 | | |
| m,p-Xylene | 42.5 | | μg/l | | 40.0 | | 106 | 70-130 | | |
| o-Xylene | 20.7 | | μg/l | | 20.0 | | 103 | 70-130 | | |
| Surrogate: 4-Bromofluorobenzene | 30.2 | | μg/l | | 30.0 | | 101 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.2 | | μg/l | | 30.0 | | 101 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 Surrogate: Dibromofluoromethane | 31.1 31.4 | | μg/l μg/l | | 30.0 30.0 | | 104 105 | 70-130 70-130 | | |
| - | 9 1.7 | | P9'' | | 00.0 | | ,00 | , 0 100 | | |
| <u>_CS Dup (8111077-BSD1)</u> | | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | 00.0 | | | | 0.5.5 | | 4 | 70 / 00 | 4- | |
| Acetone | 23.2 | | μg/l | | 20.0 | | 116 | 70-130 | 17 | 30 |

| A 1.77 | D I FI | XX :: | *DD1 | Spike | Source | A/DEC | %REC | nnn | RPD |
|---|--------------------|----------|------|-------|--------|-------|--------|--------|-------|
| Analyte(s) | Result Fl | ag Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111077 - SW846 5030 Water M | IS | | | | | | | | |
| LCS Dup (8111077-BSD1) | | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | |
| Benzene | 23.9 | μg/l | | 20.0 | | 119 | 70-130 | 10 | 30 |
| Bromodichloromethane | 21.8 | μg/l | | 20.0 | | 109 | 35-155 | 9 | 30 |
| Bromoform | 20.7 | μg/l | | 20.0 | | 103 | 45-169 | 10 | 30 |
| Bromomethane | 13.2 | μg/l | | 20.0 | | 66 | 1-242 | 21 | 30 |
| 2-Butanone (MEK) | 23.9 | μg/l | | 20.0 | | 119 | 70-130 | 24 | 30 |
| Carbon tetrachloride | 21.6 | μg/l | | 20.0 | | 108 | 70-140 | 6 | 30 |
| Chlorobenzene | 23.3 | μg/l | | 20.0 | | 116 | 70-130 | 7 | 30 |
| Chloroethane | 25.1 | μg/l | | 20.0 | | 125 | 14-230 | 2 | 30 |
| Chloroform | 24.0 | μg/l | | 20.0 | | 120 | 51-138 | 8 | 30 |
| Chloromethane | 20.3 | μg/l | | 20.0 | | 102 | 1-273 | 4 | 30 |
| Dibromochloromethane | 22.1 | μg/l | | 20.0 | | 110 | 53-149 | 9 | 30 |
| 1,2-Dichlorobenzene | 23.1 | μg/l | | 20.0 | | 115 | 18-190 | 5 | 30 |
| 1,3-Dichlorobenzene | 23.0 | μg/l | | 20.0 | | 115 | 59-156 | 6 | 30 |
| 1,4-Dichlorobenzene | 22.3 | μg/l | | 20.0 | | 112 | 18-190 | 5 | 30 |
| 1,1-Dichloroethane | 23.2 | μg/l | | 20.0 | | 116 | 59-155 | 6 | 30 |
| 1,2-Dichloroethane | 24.0 | μg/l | | 20.0 | | 120 | 49-155 | 11 | 30 |
| 1,1-Dichloroethene | 22.8 | μg/l | | 20.0 | | 114 | 70-130 | 8 | 30 |
| cis-1,2-Dichloroethene | 22.5 | μg/l | | 20.0 | | 113 | 70-130 | 9 | 30 |
| trans-1,2-Dichloroethene | 24.3 | μg/l | | 20.0 | | 122 | 54-156 | 8 | 30 |
| 1,2-Dichloropropane | 23.3 | μg/l | | 20.0 | | 117 | 1-210 | 8 | 30 |
| cis-1,3-Dichloropropene | 21.9 | μg/l | | 20.0 | | 109 | 1-227 | 7 | 30 |
| trans-1,3-Dichloropropene | 22.2 | μg/l | | 20.0 | | 111 | 17-183 | 11 | 30 |
| Ethylbenzene | 22.2 | μg/l | | 20.0 | | 111 | 37-162 | 5 | 30 |
| 2-Hexanone (MBK) | 22.8 | μg/l | | 20.0 | | 114 | 70-130 | 13 | 30 |
| Methyl tert-butyl ether | 24.8 | μg/l | | 20.0 | | 124 | 70-130 | 12 | 30 |
| 4-Methyl-2-pentanone (MIBK) | 22.3 | μg/l | | 20.0 | | 111 | 70-130 | 17 | 30 |
| Methylene chloride | 24.4 | μg/l | | 20.0 | | 122 | 1-221 | 9 | 30 |
| Styrene | 21.9 | μg/l | | 20.0 | | 109 | 70-130 | 6 | 30 |
| 1,1,2,2-Tetrachloroethane | 23.2 | μg/l | | 20.0 | | 116 | 46-157 | 11 | 30 |
| Tetrachloroethene | 22.4 | μg/l | | 20.0 | | 112 | 64-148 | 6 | 30 |
| Toluene | 22.8 | μg/l | | 20.0 | | 114 | 70-130 | 8 | 30 |
| 1,1,1-Trichloroethane | 22.7 | μg/l | | 20.0 | | 113 | 52-162 | 6 | 30 |
| 1,1,2-Trichloroethane | 23.0 | μg/l | | 20.0 | | 115 | 52-150 | 13 | 30 |
| Trichloroethene | 22.8 | μg/l | | 20.0 | | 114 | 71-157 | 8 | 30 |
| Trichlorofluoromethane (Freon 11) | 24.1 | μg/l | | 20.0 | | 120 | 17-181 | 4 | 30 |
| Vinyl chloride | 22.6 | μg/l | | 20.0 | | 113 | 1-251 | 5 | 30 |
| m,p-Xylene | 45.4 | μg/l | | 40.0 | | 113 | 70-130 | 7 | 30 |
| o-Xylene | 22.2 | μg/l | | 20.0 | | 111 | 70-130 | , 7 | 30 |
| Surrogate: 4-Bromofluorobenzene | 30.6 | μg/l | | 30.0 | | 102 | 70-130 | • | |
| Surrogate: 4 Diomondoloschizene Surrogate: Toluene-d8 | 29.8 | μg/l | | 30.0 | | 99 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 31.4 | μg/l | | 30.0 | | 105 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 31.7 | μg/l | | 30.0 | | 106 | 70-130 | | |
| Matrix Spike (8111077-MS1) | Source: SA87315-01 | | | | | | | | |
| Prepared & Analyzed: 14-Nov-08 | | | | | | | | | |
| Benzene | 17.9 | μg/l | | 20.0 | BRL | 89 | 70-130 | | |
| Bromodichloromethane | 19.6 | μg/l | | 20.0 | BRL | 98 | 35-155 | | |
| Bromoform | 17.9 | μg/l | | 20.0 | BRL | 89 | 45-169 | | |
| Bromomethane | 7.3 | μg/l | | 20.0 | BRL | 37 | 1-242 | | |
| Carbon tetrachloride | 18.8 | μg/l | | 20.0 | BRL | 94 | 70-140 | | |
| Chlorobenzene | 20.5 | μg/l | | 20.0 | BRL | 103 | 70-130 | | |

| D 1: EI | ** * | de De la companya de la companya de la companya de la companya de la companya de la companya de la companya de | Spike | Source | 0/PEG | %REC | nnn | RPD |
|--|---|--|--|---|---|--|---------------------------------|---|
| Result Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| · MS | | | | | | | | |
| Source: SA87315-01 | | | | | | | | |
| | | | | | | | | |
| 14.6 | μg/l | | 20.0 | BRL | 73 | 14-230 | | |
| 21.3 | μg/l | | 20.0 | BRL | 106 | 51-138 | | |
| 9.6 | μg/l | | 20.0 | BRL | 48 | 1-273 | | |
| 19.5 | μg/l | | 20.0 | BRL | 97 | 53-149 | | |
| 21.3 | μg/l | | 20.0 | BRL | 106 | 18-190 | | |
| 22.0 | μg/l | | 20.0 | BRL | 110 | 59-156 | | |
| 20.8 | μg/l | | 20.0 | BRL | 104 | 18-190 | | |
| 20.1 | μg/l | | 20.0 | BRL | 100 | 59-155 | | |
| 18.8 | μg/l | | 20.0 | BRL | 94 | 49-155 | | |
| 15.3 | μg/l | | 20.0 | BRL | 76 | 70-130 | | |
| 15.3 | μg/l | | 20.0 | BRL | 77 | 54-156 | | |
| 20.0 | μg/l | | 20.0 | BRL | 100 | 1-210 | | |
| 18.6 | μg/l | | 20.0 | BRL | 93 | 1-227 | | |
| 18.4 | μg/l | | 20.0 | BRL | 92 | 17-183 | | |
| 19.7 | μg/l | | 20.0 | BRL | 99 | 37-162 | | |
| 17.8 | μg/l | | 20.0 | BRL | 89 | 1-221 | | |
| 20.3 | μg/l | | 20.0 | BRL | 102 | 46-157 | | |
| 18.1 | μg/l | | 20.0 | BRL | 91 | 64-148 | | |
| 18.2 | μg/l | | 20.0 | BRL | 91 | 70-130 | | |
| 20.7 | | | 20.0 | BRL | 104 | 52-162 | | |
| 19.7 | | | 20.0 | BRL | 98 | 52-150 | | |
| 18.2 | | | 20.0 | BRL | 91 | 71-157 | | |
| 18.2 | | | 20.0 | BRL | 91 | 17-181 | | |
| 12.7 | | | 20.0 | BRL | 63 | 1-251 | | |
| 30.6 | | | 30.0 | | 102 | 70-130 | | |
| 29.6 | μg/l | | 30.0 | | 99 | 70-130 | | |
| | μg/l | | | | 103 | | | |
| 37.6 | μg/i | | 30.0 | | 105 | 70-130 | | |
| Source: SA87315-01 | | | | | | | | |
| 18.6 | μg/l | | 20.0 | BRL | 93 | 70-130 | 4 | 30 |
| 20.7 | | | 20.0 | BRL | 104 | 35-155 | 6 | 30 |
| 19.6 | μg/l | | 20.0 | BRL | 98 | 45-169 | 9 | 30 |
| 7.1 | | | 20.0 | BRL | 36 | 1-242 | 3 | 30 |
| 19.1 | | | 20.0 | BRL | 96 | 70-140 | 2 | 30 |
| 21.0 | | | 20.0 | BRL | 105 | 70-130 | 3 | 30 |
| 14.5 | | | 20.0 | BRL | 72 | 14-230 | 0.6 | 30 |
| | μg/l | | 20.0 | BRL | 110 | 51-138 | 3 | 30 |
| 21.9 | | | | | | 1-273 | 4 | 30 |
| 21.9 10.0 | μg/l | | 20.0 | BRL | 50 | 1-213 | - | |
| 10.0 | | | | | | | | |
| | μg/l | | 20.0 20.0 20.0 | BRL BRL BRL | 50 102 110 | 53-149 18-190 | 5 | 30 30 |
| 10.0 20.4 | μg/l μg/l | | 20.0 | BRL | 102 | 53-149 | 5 | 30 |
| 10.0 20.4 22.0 | μg/l μg/l μg/l | | 20.0 20.0 | BRL BRL | 102 110 | 53-149 18-190 | 5 3 | 30 30 |
| 10.0 20.4 22.0 22.8 | µg/I µg/I µg/I µg/I | | 20.0 20.0 20.0 | BRL BRL BRL | 102 110 114 | 53-149 18-190 59-156 | 5 3 4 | 30 30 30 |
| 10.0 20.4 22.0 22.8 21.4 21.0 | hā\l hā\l hā\l | | 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL | 102 110 114 107 105 | 53-149 18-190 59-156 18-190 59-155 | 5 3 4 3 4 | 30 30 30 30 30 |
| 10.0 20.4 22.0 22.8 21.4 21.0 19.8 | hā\I hā\I hā\I hā\I | | 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL | 102 110 114 107 105 99 | 53-149 18-190 59-156 18-190 59-155 49-155 | 5 3 4 3 4 5 | 30 30 30 30 30 30 |
| 10.0 20.4 22.0 22.8 21.4 21.0 19.8 15.6 | hā\I hā\I hā\I hā\I hā\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL | 102 110 114 107 105 99 78 | 53-149 18-190 59-156 18-190 59-155 49-155 70-130 | 5 3 4 3 4 5 | 30 30 30 30 30 30 30 |
| 10.0 20.4 22.0 22.8 21.4 21.0 19.8 15.6 | ha\] ha\] ha\] ha\] ha\] ha\] | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL | 102 110 114 107 105 99 78 79 | 53-149 18-190 59-156 18-190 59-155 49-155 70-130 54-156 | 5 3 4 3 4 5 2 | 30 30 30 30 30 30 30 30 |
| 10.0 20.4 22.0 22.8 21.4 21.0 19.8 15.6 | hā\I hā\I hā\I hā\I hā\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL | 102 110 114 107 105 99 78 | 53-149 18-190 59-156 18-190 59-155 49-155 70-130 | 5 3 4 3 4 5 | 30 30 30 30 30 30 30 |
| | \$\text{Source: SA87315-01}\$ 14.6 21.3 9.6 19.5 21.3 22.0 20.8 20.1 18.8 15.3 15.3 20.0 18.6 18.4 19.7 17.8 20.3 18.1 18.2 20.7 19.7 18.2 18.2 12.7 30.6 29.6 30.8 31.6 \$\text{Source: SA87315-01}\$ 18.6 20.7 19.6 7.1 19.1 21.0 14.5 | MS Source: SA87315-01 14.6 | MS Source: SA87315-01 14.6 | NS Source: SA87315-01 | Result Flag Units *RDL Level Result | Result Flag Units *RDL Level Result %REC | MS Source: SA87315-01 14.6 | Result Flag Units *RDL Level Result *REC Limits RPD |

| Result Flag e: SA87315-01 20.3 18.5 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 BRL BRL | Units µg/l µg/l µg/l µg/l µg/l µg/l µg/l µg/ | *RDL | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL BRL | %REC 102 92 107 95 93 108 106 92 92 63 101 101 105 105 | 37-162 1-221 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | RPD 3 4 5 4 2 4 8 1 2 1 | 30 30 30 30 30 30 30 30 30 30 30 |
|---|--|-------------|---|--|--|--|---------------------------------|--|
| 20.3 18.5 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL | 92 107 95 93 108 106 92 92 63 101 101 | 1-221 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 5 4 2 4 8 1 | 30 30 30 30 30 30 30 30 |
| 20.3 18.5 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL | 92 107 95 93 108 106 92 92 63 101 101 | 1-221 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 5 4 2 4 8 1 | 30 30 30 30 30 30 30 30 |
| 18.5 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL | 92 107 95 93 108 106 92 92 63 101 101 | 1-221 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 5 4 2 4 8 1 | 30 30 30 30 30 30 30 30 |
| 18.5 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL BRL | 92 107 95 93 108 106 92 92 63 101 101 | 1-221 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 5 4 2 4 8 1 | 30 30 30 30 30 30 30 30 |
| 21.4 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\l hâ\l hâ\l hâ\l hâ\l hâ\l hâ\l | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 20.0 | BRL BRL BRL BRL BRL BRL | 107 95 93 108 106 92 92 63 101 101 105 | 46-157 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 5 4 2 4 8 1 | 30 30 30 30 30 30 30 |
| 19.0 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 20.0 30.0 3 | BRL BRL BRL BRL BRL BRL | 95 93 108 106 92 92 63 101 101 105 | 64-148 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 2 4 8 1 2 | 30 30 30 30 30 30 |
| 18.7 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 20.0 30.0 30.0 | BRL BRL BRL BRL BRL | 93 108 106 92 92 63 101 101 105 | 70-130 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 2 4 8 1 2 | 30 30 30 30 30 |
| 21.5 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 20.0 30.0 30.0 30.0 | BRL BRL BRL BRL | 108 106 92 92 63 101 101 105 | 52-162 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 4 8 1 2 | 30 30 30 30 |
| 21.3 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 20.0 30.0 30.0 30.0 | BRL BRL BRL | 106 92 92 63 101 101 105 | 52-150 71-157 17-181 1-251 70-130 70-130 70-130 | 8 1 2 | 30 30 30 |
| 18.4 18.5 12.5 30.4 30.2 31.5 31.6 | hâ\I hâ\I hâ\I hâ\I hâ\I | | 20.0 20.0 20.0 30.0 30.0 30.0 | BRL BRL | 92 92 63 101 101 105 | 71-157 17-181 1-251 70-130 70-130 70-130 | 1 2 | 30 30 |
| 18.5 12.5 30.4 30.2 31.5 31.6 BRL | hâ\l hâ\l hâ\l hâ\l | | 20.0 20.0 30.0 30.0 30.0 | BRL | 92 63 101 101 105 | 17-181 1-251 70-130 70-130 70-130 | 2 | 30 |
| 12.5 30.4 30.2 31.5 31.6 BRL BRL | hâ\I hâ\I hâ\I hâ\I | | 20.0 30.0 30.0 30.0 | | 63 101 101 105 | 1-251 70-130 70-130 70-130 | | |
| 30.4 30.2 31.5 31.6 BRL | hâ\I hâ\I hâ\I hâ\I | | 30.0 30.0 30.0 | BRL | 101 101 105 | 70-130 70-130 70-130 | 1 | 30 |
| 30.2 31.5 31.6 BRL BRL | hâ\I hâ\I hâ\I hâ\I | | 30.0 30.0 | | 101 105 | 70-130 70-130 | | |
| 31.5 31.6 BRL BRL | µg/I µg/I µg/I | | 30.0 | | 105 | 70-130 | | |
| 31.6 BRL BRL | μg/l μg/l | | | | | | | |
| BRL BRL | μg/l | | 30.0 | | 105 | 70-130 | | |
| BRL | | | | | | | | |
| BRL | | | | | | | | |
| BRL | | | | | | | | |
| BRL | | 20.0 | | | | | | |
| | | | | | | | | |
| חחו | μg/l | 1.0 | | | | | | |
| BRL | μg/l | 1.0 | | | | | | |
| BRL | μg/l | 1.0 | | | | | | |
| BRL | μg/l | 2.0 | | | | | | |
| BRL BRL | μg/l | 10.0 1.0 | | | | | | |
| | μg/l | | | | | | | |
| BRL | μg/l | 1.0 | | | | | | |
| BRL | μg/l | 2.0 | | | | | | |
| BRL | μg/l | 1.0 | | | | | | |
| BRL | μg/l | 2.0 | | | | | | |
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| | μg/l | | | | | | | |
| | μg/l | | | | | | | |
| BRL | μg/l | | | | | | | |
| | μg/l | | | | | | | |
| | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | BRL | BRL | BRL μg/l 1.0 | BRL | BRL | BRL | BRL |

| | | | | | Spike | Source | | %REC | | RPD |
|--|--------------|------|--------------|------|--------------|--------|-----------|------------------|-----|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111194 - SW846 5030 Water MS | | | | | | | | | | |
| Blank (8111194-BLK1) | | | | | | | | | | |
| Prepared & Analyzed: 17-Nov-08 | | | | | | | | | | |
| 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| Trichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | | | | | | |
| Vinyl chloride | BRL | | μg/l | 1.0 | | | | | | |
| m,p-Xylene | BRL | | μg/l | 2.0 | | | | | | |
| o-Xylene | BRL | | μg/l | 1.0 | | | | | | |
| Surrogate: 4-Bromofluorobenzene | 25.9 | | μg/l | | 30.0 | | 86 | 70-130 | | |
| Surrogate: Toluene-d8 Surrogate: 1,2-Dichloroethane-d4 | 29.6 33.2 | | μg/l μg/l | | 30.0 30.0 | | 99 111 | 70-130 70-130 | | |
| Surrogate: Dibromofluoromethane | 34.4 | | μg/l | | 30.0 | | 115 | 70-130 | | |
| LCS (8111194-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 17-Nov-08 | | | | | | | | | | |
| Acetone | 18.5 | | μg/l | | 20.0 | | 92 | 70-130 | | |
| Benzene | 20.2 | | μg/l | | 20.0 | | 101 | 70-130 | | |
| Bromodichloromethane | 24.1 | | μg/l | | 20.0 | | 121 | 35-155 | | |
| Bromoform | 18.7 | | μg/l | | 20.0 | | 93 | 45-169 | | |
| Bromomethane | 19.8 | | μg/l | | 20.0 | | 99 | 1-242 | | |
| 2-Butanone (MEK) | 20.0 | | μg/l | | 20.0 | | 100 | 70-130 | | |
| Carbon tetrachloride | 20.7 | | μg/l | | 20.0 | | 103 | 70-140 | | |
| Chlorobenzene | 20.3 | | μg/l | | 20.0 | | 102 | 70-130 | | |
| Chloroethane | 20.6 | | μg/l | | 20.0 | | 103 | 14-230 | | |
| Chloroform | 22.8 | | μg/l | | 20.0 | | 114 | 51-138 | | |
| Chloromethane | 18.6 | | μg/l | | 20.0 | | 93 | 1-273 | | |
| Dibromochloromethane | 20.0 | | μg/l | | 20.0 | | 100 | 53-149 | | |
| 1,2-Dichlorobenzene | 18.8 | | μg/l | | 20.0 | | 94 | 18-190 | | |
| 1,3-Dichlorobenzene | 22.0 | | μg/l | | 20.0 | | 110 | 59-156 | | |
| 1,4-Dichlorobenzene | 18.9 | | μg/l | | 20.0 | | 94 | 18-190 | | |
| 1,1-Dichloroethane | 20.4 | | μg/l | | 20.0 | | 102 | 59-155 | | |
| 1,2-Dichloroethane | 20.4 | | μg/l | | 20.0 | | 102 | 49-155 | | |
| 1,1-Dichloroethene | 17.5 | | μg/l | | 20.0 | | 87 | 70-130 | | |
| cis-1,2-Dichloroethene | 19.2 | | μg/l | | 20.0 | | 96 | 70-130 | | |
| trans-1,2-Dichloroethene | 19.6 | | μg/l | | 20.0 | | 98 | 54-156 | | |
| 1,2-Dichloropropane | 20.7 | | μg/l | | 20.0 | | 103 | 1-210 | | |
| cis-1,3-Dichloropropene | 18.4 | | μg/l | | 20.0 | | 92 | 1-227 | | |
| trans-1,3-Dichloropropene | 19.0 | | μg/l | | 20.0 | | 95 | 17-183 | | |
| Ethylbenzene | 20.1 | | μg/l | | 20.0 | | 100 | 37-162 | | |
| 2-Hexanone (MBK) | 14.5 | | μg/l | | 20.0 | | 73 | 70-130 | | |
| Methyl tert-butyl ether | 21.2 | | μg/l | | 20.0 | | 106 | 70-130 | | |
| 4-Methyl-2-pentanone (MIBK) | 19.4 | | μg/l | | 20.0 | | 97 | 70-130 | | |
| Methylene chloride | 21.6 | | μg/l | | 20.0 | | 108 | 1-221 | | |
| Styrene | 18.3 | | μg/l | | 20.0 | | 92 | 70-130 | | |
| 1,1,2,2-Tetrachloroethane | 21.3 | | μg/l | | 20.0 | | 106 | 46-157 | | |
| Tetrachloroethene | 16.8 | | μg/l | | 20.0 | | 84 | 64-148 | | |
| Toluene | 19.8 | | μg/l | | 20.0 | | 99 | 70-130 | | |
| 1,1,1-Trichloroethane | 18.7 | | μg/l | | 20.0 | | 94 | 52-162 | | |
| 1,1,2-Trichloroethane | 21.4 | | μg/l | | 20.0 | | 107 | 52-150 | | |
| Trichloroethene | 19.6 | | μg/l | | 20.0 | | 98 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 17.0 | | μg/l | | 20.0 | | 85 | 17-181 | | |
| Vinyl chloride | 19.7 | | μg/l | | 20.0 | | 99 | 1-251 | | |

| | | Source | | %REC | | RPD |
|------|--------------|----------------------|------------------------------|---|---|---|
| *RDL | Level | Result | %REC | Limits | RPD | Limit |
| | | | | | | |
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| | | | | | | |
| | 40.0 | | 105 | 70-130 | | |
| | 20.0 | | 105 | 70-130 | | |
| | 30.0 | | 106 | 70-130 | | |
| | 30.0 30.0 | | 101 103 | 70-130 70-130 | | |
| | 30.0 | | 103 | 70-130 70-130 | | |
| | | | | | | |
| | | | | | | |
| | 20.0 | | 95 | 70-130 | 3 | 30 |
| | 20.0 | | 111 | 70-130 | 9 | 30 |
| | 20.0 | | 131 | 35-155 | 9 | 30 |
| | 20.0 | | 96 | 45-169 | 3 | 30 |
| | 20.0 | | 120 | 1-242 | 19 | 30 |
| | 20.0 | | 98 | 70-130 | 2 | 30 |
| | 20.0 | | 127 | 70-140 | 20 | 30 |
| | 20.0 | | 112 | 70-130 | 10 | 30 |
| | 20.0 | | 119 | 14-230 | 14 | 30 |
| | 20.0 | | 126 | 51-138 | 10 | 30 |
| | 20.0 | | 105 | 1-273 | 12 | 30 |
| | 20.0 | | 105 | 53-149 | 5 | 30 |
| | 20.0 | | 104 | 18-190 | 10 | 30 |
| | 20.0 | | 121 | 59-156 | 10 | 30 |
| | 20.0 | | 105 | 18-190 | 11 | 30 |
| | 20.0 | | 113 | 59-155 | 10 | 30 |
| | 20.0 | | 109 | 49-155 | 6 | 30 |
| | 20.0 | | 102 | 70-130 | 15 | 30 |
| | 20.0 | | 106 | 70-130 | 9 | 30 |
| | 20.0 | | 111 | 54-156 | 12 | 30 |
| | 20.0 | | 110 | 1-210 | 7 | 30 |
| | 20.0 | | 95 | 1-227 | 4 | 30 |
| | 20.0 | | 96 | 17-183 | 8.0 | 30 |
| | 20.0 | | 115 | 37-162 | 13 | 30 |
| | 20.0 | | 68 | 70-130 | 6 | 30 |
| | 20.0 | | 106 | 70-130 | 0.5 | 30 |
| | 20.0 | | 90 | 70-130 | 8 | 30 |
| | 20.0 | | 119 | 1-221 | 10 | 30 |
| | 20.0 | | 97 | 70-130 | 6 | 30 |
| | 20.0 | | 110 | 46-157 | 3 | 30 |
| | 20.0 | | 100 | 64-148 | 17 | 30 |
| | 20.0 | | 111 | 70-130 | 11 | 30 |
| | 20.0 | | 108 | 52-162 | 15 | 30 |
| | 20.0 | | 109 | 52-150 | 2 | 30 |
| | 20.0 | | 110 | 71-157 | 12 | 30 |
| | 20.0 | | 108 | 17-181 | 24 | 30 |
| | 20.0 | | 147 | 1-251 | 39 | 30 |
| | 40.0 | | 119 | 70-130 | 13 | 30 |
| | 20.0 | | 119 | 70-130 | 12 | 30 |
| | 30.0 | | 108 | 70-130 | | - |
| | 30.0 | | 101 | 70-130 | | |
| | | 40.0 20.0 30.0 | 40.0 20.0 30.0 30.0 | 40.0 119 20.0 119 30.0 108 30.0 101 | 40.0 119 70-130 20.0 119 70-130 30.0 108 70-130 30.0 101 70-130 | 40.0 119 70-130 13 20.0 119 70-130 12 30.0 108 70-130 30.0 101 70-130 |

| Analyto(a) | Result | Flag | Unita | *RDL | Spike | Source | %REC | %REC | RPD | RPD Limit |
|---|-----------------|------|-------|------|-------|--------|-------|--------|-----|--------------|
| Analyte(s) | Result | гіад | Units | ·KDL | Level | Result | 70KEC | Limits | KPD | Limit |
| Batch 8111194 - SW846 5030 Water | MS | | | | | | | | | |
| LCS Dup (8111194-BSD1) Prepared & Analyzed: 17-Nov-08 | | | | | | | | | | |
| Surrogate: Dibromofluoromethane | 32.1 | | μg/l | | 30.0 | | 107 | 70-130 | | |
| Matrix Spike (8111194-MS1) | Source: SA87191 | -03 | | | | | | | | |
| Prepared & Analyzed: 17-Nov-08 | Cource. CAO? 13 | -00 | | | | | | | | |
| Benzene | 12.8 | QM7 | μg/l | | 20.0 | BRL | 64 | 70-130 | | |
| Bromodichloromethane | 24.2 | | μg/l | | 20.0 | BRL | 121 | 35-155 | | |
| Bromoform | 19.6 | | μg/l | | 20.0 | BRL | 98 | 45-169 | | |
| Bromomethane | 6.1 | | μg/l | | 20.0 | BRL | 30 | 1-242 | | |
| Carbon tetrachloride | 17.9 | | μg/l | | 20.0 | BRL | 89 | 70-140 | | |
| Chlorobenzene | 18.4 | | μg/l | | 20.0 | BRL | 92 | 70-130 | | |
| Chloroethane | 8.6 | | μg/l | | 20.0 | BRL | 43 | 14-230 | | |
| Chloroform | 21.9 | | μg/l | | 20.0 | BRL | 109 | 51-138 | | |
| Chloromethane | 4.4 | | μg/l | | 20.0 | BRL | 22 | 1-273 | | |
| Dibromochloromethane | 20.3 | | μg/l | | 20.0 | BRL | 101 | 53-149 | | |
| 1,2-Dichlorobenzene | 21.9 | | μg/l | | 20.0 | BRL | 109 | 18-190 | | |
| 1,3-Dichlorobenzene | 24.4 | | μg/l | | 20.0 | BRL | 122 | 59-156 | | |
| 1,4-Dichlorobenzene | 19.8 | | μg/l | | 20.0 | BRL | 99 | 18-190 | | |
| 1,1-Dichloroethane | 17.6 | | μg/l | | 20.0 | BRL | 88 | 59-155 | | |
| 1,2-Dichloroethane | 18.4 | | μg/l | | 20.0 | BRL | 92 | 49-155 | | |
| 1,1-Dichloroethene | 9.2 | QM7 | μg/l | | 20.0 | BRL | 46 | 70-130 | | |
| rans-1,2-Dichloroethene | 9.2 | QM7 | μg/l | | 20.0 | BRL | 46 | 54-156 | | |
| 1,2-Dichloropropane | 18.0 | | μg/l | | 20.0 | BRL | 90 | 1-210 | | |
| cis-1,3-Dichloropropene | 15.2 | | μg/l | | 20.0 | BRL | 76 | 1-227 | | |
| rans-1,3-Dichloropropene | 16.9 | | μg/l | | 20.0 | BRL | 84 | 17-183 | | |
| Ethylbenzene | 17.1 | | μg/l | | 20.0 | BRL | 86 | 37-162 | | |
| Methylene chloride | 14.4 | | μg/l | | 20.0 | BRL | 72 | 1-221 | | |
| 1,1,2,2-Tetrachloroethane | 23.7 | | μg/l | | 20.0 | BRL | 118 | 46-157 | | |
| Tetrachloroethene | 12.6 | QM7 | μg/l | | 20.0 | BRL | 63 | 64-148 | | |
| Toluene | 15.3 | | μg/l | | 20.0 | BRL | 76 | 70-130 | | |
| 1,1,1-Trichloroethane | 20.2 | | μg/l | | 20.0 | 1.3 | 95 | 52-162 | | |
| 1,1,2-Trichloroethane | 22.3 | | μg/l | | 20.0 | BRL | 111 | 52-150 | | |
| Trichloroethene | 14.1 | QM7 | μg/l | | 20.0 | BRL | 70 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 12.4 | | μg/l | | 20.0 | BRL | 62 | 17-181 | | |
| Vinyl chloride | 5.9 | | μg/l | | 20.0 | BRL | 29 | 1-251 | | |
| Surrogate: 4-Bromofluorobenzene | 33.5 | | μg/l | | 30.0 | | 112 | 70-130 | | |
| Surrogate: Toluene-d8 | 31.0 | | μg/l | | 30.0 | | 103 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 34.5 | | μg/l | | 30.0 | | 115 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 33.6 | | μg/l | | 30.0 | | 112 | 70-130 | | |
| Matrix Spike Dup (8111194-MSD1) | Source: SA87191 | -03 | | | | | | | | |
| Prepared & Analyzed: 17-Nov-08 | | | | | | | | | | |
| Benzene | 12.1 | QM7 | μg/l | | 20.0 | BRL | 61 | 70-130 | 6 | 30 |
| Bromodichloromethane | 21.7 | | μg/l | | 20.0 | BRL | 109 | 35-155 | 11 | 30 |
| Bromoform | 17.6 | | μg/l | | 20.0 | BRL | 88 | 45-169 | 11 | 30 |
| Bromomethane | 6.3 | | μg/l | | 20.0 | BRL | 31 | 1-242 | 3 | 30 |
| Carbon tetrachloride | 15.6 | | μg/l | | 20.0 | BRL | 78 | 70-140 | 14 | 30 |
| Chlorobenzene | 17.1 | | μg/l | | 20.0 | BRL | 85 | 70-130 | 7 | 30 |
| Chloroethane | 8.1 | | μg/l | | 20.0 | BRL | 41 | 14-230 | 6 | 30 |
| Chloroform | 19.8 | | μg/l | | 20.0 | BRL | 99 | 51-138 | 10 | 30 |
| Chloromethane | 4.1 | | μg/l | | 20.0 | BRL | 21 | 1-273 | 7 | 30 |
| Dibromochloromethane | 18.4 | | μg/l | | 20.0 | BRL | 92 | 53-149 | 10 | 30 |

| | | | | | Spike | Source | | %REC | | RPD |
|--|---|------------|--|--|--------------|--------|------------|------------------|---------|------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limi |
| Batch 8111194 - SW846 5030 Water | r MS | | | | | | | | | |
| Matrix Spike Dup (8111194-MSD1) | Source: SA8719 | 1-03 | | | | | | | | |
| Prepared & Analyzed: 17-Nov-08 | | | | | | | | | | |
| 1,2-Dichlorobenzene | 20.2 | | μg/l | | 20.0 | BRL | 101 | 18-190 | 8 | 30 |
| 1,3-Dichlorobenzene | 22.1 | | μg/l | | 20.0 | BRL | 111 | 59-156 | 10 | 30 |
| 1,4-Dichlorobenzene | 18.6 | | μg/l | | 20.0 | BRL | 93 | 18-190 | 6 | 30 |
| 1,1-Dichloroethane | 16.2 | | μg/l | | 20.0 | BRL | 81 | 59-155 | 9 | 30 |
| 1,2-Dichloroethane | 16.5 | | μg/l | | 20.0 | BRL | 82 | 49-155 | 11 | 30 |
| 1,1-Dichloroethene | 8.4 | QM7 | μg/l | | 20.0 | BRL | 42 | 70-130 | 8 | 30 |
| trans-1,2-Dichloroethene | 8.7 | QM7 | μg/l | | 20.0 | BRL | 44 | 54-156 | 5 | 30 |
| 1,2-Dichloropropane | 16.9 | | μg/l | | 20.0 | BRL | 84 | 1-210 | 7 | 30 |
| cis-1,3-Dichloropropene | 14.3 | | μg/l | | 20.0 | BRL | 72 | 1-227 | 6 | 30 |
| trans-1,3-Dichloropropene | 15.5 | | | | 20.0 | BRL | 77 | 17-183 | 9 | 30 |
| Ethylbenzene | 15.9 | | μg/l | | 20.0 | BRL | 80 | 37-162 | 7 | 30 |
| Methylene chloride | 13.1 | | μg/l | | 20.0 | BRL | 65 | 1-221 | 10 | 30 |
| 1,1,2,2-Tetrachloroethane | 21.8 | | μg/l | | 20.0 | BRL | 109 | 46-157 | 8 | 30 |
| Tetrachloroethene | | QM7 | µg/l | | | BRL | | | 8 7 | 30 |
| Toluene | 11.8 13.9 | QM7 QC1 | μg/l | | 20.0 20.0 | BRL | 59 69 | 64-148 70-130 | 7 10 | 30 |
| | | QUI | µg/l | | | | | | | |
| 1,1,1-Trichloroethane | 18.1 | | µg/l | | 20.0 | 1.3 | 84 | 52-162 | 12 | 30 |
| 1,1,2-Trichloroethane | 20.2 | 01.17 | μg/l | | 20.0 | BRL | 101 | 52-150 | 10 | 30 |
| Trichloroethene | 13.2 | QM7 | μg/l | | 20.0 | BRL | 66 | 71-157 | 6 | 30 |
| Trichlorofluoromethane (Freon 11) | 11.0 | | µg/l | | 20.0 | BRL | 55 | 17-181 | 12 | 30 |
| Vinyl chloride | 5.2 | | μg/l | | 20.0 | BRL | 26 | 1-251 | 13 | 30 |
| Surrogate: 4-Bromofluorobenzene Surrogate: Toluene-d8 | 32.2 30.2 | | µg/l µg/l | | 30.0 30.0 | | 107 101 | 70-130 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 32.8 | | μg/l | | 30.0 | | 109 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 32.8 | | μg/l | | 00.0 | | 400 | 70-130 | | |
| = | | | μул | | 30.0 | | 109 | 70-730 | | |
| Batch 8111291 - SW846 5030 Water | | | рул | | 30.0 | | 109 | 70-730 | | |
| Batch 8111291 - SW846 5030 Water | | | μул | | 30.0 | | 109 | 70-730 | | |
| | | | μg/i | | 30.0 | | 109 | 70-730 | | |
| Blank (8111291-BLK1) | | | µg/I | 20.0 | 30.0 | | 109 | 70-730 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 | r MS | | | 20.0 1.0 | 30.0 | | 109 | 70-730 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone | r MS BRL | | μg/l | | 30.0 | | 109 | 70-730 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene | r MS BRL BRL | | µg/l µg/l | 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane | BRL BRL BRL | | µg/l µg/l µg/l | 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform | BRL BRL BRL BRL BRL | | hā\] hā\] hā\] | 1.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane | BRL BRL BRL BRL BRL BRL | | hā\] hā\] hā\] | 1.0 1.0 1.0 2.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) | BRL BRL BRL BRL BRL BRL BRL | | hā\] hā\] hā\] hā\] | 1.0 1.0 1.0 2.0 10.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride | BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene | BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroethane Chloroform | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroform Chloromethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 2.0 10.0 1.0 1.0 2.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroform Chloromethane Dibromochloromethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 2.0 10.0 1.0 1.0 2.0 1.0 2.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroethane Chloroform Chloromethane Dibromochloromethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroethane Chloroform Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,3-Dichlorobenzene | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | hall hall hall hall hall hall hall hall | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Blank (8111291-BLK1) Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroethane Chloroform Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chloroform Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,1-Dichloroethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chlorothane Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,4-Dichlorobenzene 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | hall hall hall hall hall hall hall hall | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |
| Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromoform Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chlorotemane Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,4-Dichlorobenzene 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethene 1,1-Dichloroethene | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ ha\/ | 1.0 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1 | 30.0 | | 109 | 70-130 | | |
| Prepared & Analyzed: 18-Nov-08 Acetone Benzene Bromodichloromethane Bromomethane 2-Butanone (MEK) Carbon tetrachloride Chlorobenzene Chlorothane Chloromethane Dibromochloromethane 1,2-Dichlorobenzene 1,4-Dichlorobenzene 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane | BRL BRL BRL BRL BRL BRL BRL BRL BRL BRL | | hall hall hall hall hall hall hall hall | 1.0 1.0 1.0 2.0 10.0 1.0 2.0 1.0 2.0 1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 | 30.0 | | 109 | 70-130 | | |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|--|--------------|------|--------------|------|----------------|------------------|-----------|------------------|-----|--------------|
| Batch 8111291 - SW846 5030 Water MS | | | | | | | | | | |
| Blank (8111291-BLK1) | | | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| trans-1,3-Dichloropropene | BRL | | μg/l | 1.0 | | | | | | |
| Ethylbenzene | BRL | | μg/l | 1.0 | | | | | | |
| 2-Hexanone (MBK) | BRL | | μg/l | 10.0 | | | | | | |
| Methyl tert-butyl ether | BRL | | μg/l | 1.0 | | | | | | |
| 4-Methyl-2-pentanone (MIBK) | BRL | | μg/l | 10.0 | | | | | | |
| Methylene chloride | BRL | | μg/l | 10.0 | | | | | | |
| Styrene | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,2,2-Tetrachloroethane | BRL | | μg/l | 1.0 | | | | | | |
| Tetrachloroethene | BRL | | μg/l | 1.0 | | | | | | |
| Toluene | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,1-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| 1,1,2-Trichloroethane | BRL | | μg/l | 1.0 | | | | | | |
| Trichloroethene | BRL | | μg/l | 1.0 | | | | | | |
| Trichlorofluoromethane (Freon 11) | BRL | | μg/l | 1.0 | | | | | | |
| Vinyl chloride | BRL | | μg/l | 1.0 | | | | | | |
| m,p-Xylene | BRL | | μg/l | 2.0 | | | | | | |
| o-Xylene | BRL | | μg/l | 1.0 | | | | | | |
| Surrogate: 4-Bromofluorobenzene | 25.1 | | μg/l | | 30.0 | | 84 | 70-130 | | |
| Surrogate: Toluene-d8 Surrogate: 1,2-Dichloroethane-d4 | 29.7 33.8 | | μg/l μg/l | | 30.0 30.0 | | 99 113 | 70-130 70-130 | | |
| Surrogate: Dibromofluoromethane | 36.4 | | μg/l | | 30.0 | | 121 | 70-130 | | |
| LCS (8111291-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| Acetone | 20.6 | | μg/l | | 20.0 | | 103 | 70-130 | | |
| Benzene | 20.6 | | μg/l | | 20.0 | | 103 | 70-130 | | |
| Bromodichloromethane | 24.1 | | μg/l | | 20.0 | | 121 | 35-155 | | |
| Bromoform | 18.3 | | μg/l | | 20.0 | | 92 | 45-169 | | |
| Bromomethane | 23.3 | | μg/l | | 20.0 | | 116 | 1-242 | | |
| 2-Butanone (MEK) | 19.0 | | μg/l | | 20.0 | | 95 | 70-130 | | |
| Carbon tetrachloride | 22.8 | | μg/l | | 20.0 | | 114 | 70-140 | | |
| Chlorobenzene | 20.8 | | μg/l | | 20.0 | | 104 | 70-130 | | |
| Chloroethane | 22.1 | | μg/l | | 20.0 | | 110 | 14-230 | | |
| Chloroform | 23.2 | | μg/l | | 20.0 | | 116 | 51-138 | | |
| Chloromethane | 19.7 | | μg/l | | 20.0 | | 99 | 1-273 | | |
| Dibromochloromethane | 19.7 | | μg/l | | 20.0 | | 98 | 53-149 | | |
| 1,2-Dichlorobenzene | 19.2 | | μg/l | | 20.0 | | 96 | 18-190 | | |
| 1,3-Dichlorobenzene | 22.7 | | μg/l | | 20.0 | | 113 | 59-156 | | |
| 1,4-Dichlorobenzene | 19.3 | | μg/l | | 20.0 | | 96 | 18-190 | | |
| 1,1-Dichloroethane | 21.0 | | μg/l | | 20.0 | | 105 | 59-155 | | |
| 1,2-Dichloroethane | 20.6 | | μg/l | | 20.0 | | 103 | 49-155 | | |
| 1,1-Dichloroethene | 19.1 | | μg/l | | 20.0 | | 96 | 70-130 | | |
| cis-1,2-Dichloroethene | 19.7 | | μg/l | | 20.0 | | 99 | 70-130 | | |
| trans-1,2-Dichloroethene | 20.7 | | μg/l | | 20.0 | | 103 | 54-156 | | |
| 1,2-Dichloropropane | 20.8 | | μg/l | | 20.0 | | 104 | 1-210 | | |
| cis-1,3-Dichloropropene | 17.8 | | μg/l | | 20.0 | | 89 | 1-227 | | |
| trans-1,3-Dichloropropene | 17.9 | | μg/l | | 20.0 | | 89 | 17-183 | | |
| Ethylbenzene | 21.1 | | μg/l | | 20.0 | | 106 | 37-162 | | |
| 2-Hexanone (MBK) | 14.0 | | μg/l | | 20.0 | | 70 | 70-130 | | |
| Methyl tert-butyl ether | 21.0 | | μg/l | | 20.0 | | 105 | 70-130 | | |
| 4-Methyl-2-pentanone (MIBK) | 18.0 | | μg/l | | 20.0 | | 90 | 70-130 | | |

| | | | | | Spike | Source | | %REC | | RPD |
|--|--------------|------|--------------|------|--------------|--------|------------|------------------|----------|----------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111291 - SW846 5030 Water MS | | | | | | | | | | |
| LCS (8111291-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| Methylene chloride | 22.5 | | μg/l | | 20.0 | | 113 | 1-221 | | |
| Styrene | 18.0 | | μg/l | | 20.0 | | 90 | 70-130 | | |
| 1,1,2,2-Tetrachloroethane | 21.1 | | μg/l | | 20.0 | | 106 | 46-157 | | |
| Tetrachloroethene | 18.1 | | μg/l | | 20.0 | | 90 | 64-148 | | |
| Toluene | 20.4 | | μg/l | | 20.0 | | 102 | 70-130 | | |
| 1,1,1-Trichloroethane | 19.9 | | μg/l | | 20.0 | | 100 | 52-162 | | |
| 1,1,2-Trichloroethane | 21.0 | | μg/l | | 20.0 | | 105 | 52-150 | | |
| Trichloroethene | 20.1 | | μg/l | | 20.0 | | 101 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 19.6 | | μg/l | | 20.0 | | 98 | 17-181 | | |
| Vinyl chloride | 23.2 | | μg/l | | 20.0 | | 116 | 1-251 | | |
| m,p-Xylene | 43.9 | | μg/l | | 40.0 | | 110 | 70-130 | | |
| o-Xylene | 22.1 | | μg/l | | 20.0 | | 111 | 70-130 | | |
| Surrogate: 4-Bromofluorobenzene | 32.0 | | μg/l | | 30.0 | | 107 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.3 | | μg/l | | 30.0 | | 101 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 Surrogate: Dibromofluoromethane | 31.3 32.1 | | μg/l μg/l | | 30.0 30.0 | | 104 107 | 70-130 70-130 | | |
| - | | | F-3·· | | | | | | | |
| LCS Dup (8111291-BSD1) | | | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | 0.4.4 | | | | 00.0 | | 407 | 70.400 | | 00 |
| Acetone | 21.4 | | μg/l | | 20.0 | | 107 | 70-130 | 4 | 30 |
| Benzene | 20.1 | | μg/l | | 20.0 | | 100 | 70-130 | 3 | 30 |
| Bromodichloromethane | 23.8 | | μg/l | | 20.0 | | 119 | 35-155 | 1 | 30 |
| Bromoform | 18.8 | | μg/l | | 20.0 | | 94 | 45-169 | 2 | 30 |
| Bromomethane | 22.6 | | μg/l | | 20.0 | | 113 | 1-242 | 3 | 30 |
| 2-Butanone (MEK) | 20.9 | | μg/l | | 20.0 | | 104 | 70-130 | 10 | 30 |
| Carbon tetrachloride | 21.8 | | μg/l | | 20.0 | | 109 | 70-140 | 4 | 30 |
| Chlorothore | 20.5 | | μg/l | | 20.0 | | 103 | 70-130 | 1 | 30 |
| Chloroethane Chloroform | 21.0 23.3 | | μg/l | | 20.0 20.0 | | 105 117 | 14-230 51-138 | 5 0.6 | 30 30 |
| | 23.3 18.6 | | µg/l | | 20.0 | | | | | |
| Chloromethane Dibromochloromethane | 19.6 | | μg/l | | | | 93 98 | 1-273 53-149 | 6 0.1 | 30 30 |
| | | | μg/l | | 20.0 | | | | | |
| 1,2-Dichlorobenzene | 19.6 | | µg/l | | 20.0 | | 98 | 18-190 | 2 | 30 |
| 1,3-Dichlorobenzene 1,4-Dichlorobenzene | 22.9 20.0 | | μg/l | | 20.0 | | 114 | 59-156 | 0.7 | 30 30 |
| 1,1-Dichloroethane | 20.6 | | μg/l | | 20.0 20.0 | | 100 103 | 18-190 59-155 | 4 | 30 |
| 1,2-Dichloroethane | 21.0 | | µg/l | | 20.0 | | 105 | 49-155 | 2 2 | 30 |
| 1,1-Dichloroethene | 18.3 | | µg/l | | 20.0 | | 92 | 70-130 | 4 | 30 |
| cis-1,2-Dichloroethene | 19.3 | | µg/l | | 20.0 | | 96 | 70-130 | 2 | 30 |
| trans-1,2-Dichloroethene | 19.9 | | µg/l | | 20.0 | | 99 | 54-156 | 4 | 30 |
| 1,2-Dichloropropane | 20.4 | | μg/l μg/l | | 20.0 | | 102 | 1-210 | 2 | 30 |
| cis-1,3-Dichloropropene | 17.7 | | μg/l | | 20.0 | | 88 | 1-217 | 0.4 | 30 |
| trans-1,3-Dichloropropene | 18.0 | | μg/l | | 20.0 | | 90 | 17-183 | 0.4 | 30 |
| Ethylbenzene | 20.1 | | | | 20.0 | | 100 | 37-162 | 5 | 30 |
| 2-Hexanone (MBK) | 15.4 | | μg/l μg/l | | 20.0 | | 77 | 70-130 | 9 | 30 |
| Methyl tert-butyl ether | 21.6 | | μg/l | | 20.0 | | 108 | 70-130 | 3 | 30 |
| 4-Methyl-2-pentanone (MIBK) | 19.2 | | μg/l | | 20.0 | | 96 | 70-130 | 3 7 | 30 |
| Methylene chloride | 22.3 | | μg/l | | 20.0 | | 111 | 1-221 | 1 | 30 |
| Styrene | 17.6 | | μg/l | | 20.0 | | 88 | 70-130 | 2 | 30 |
| 1,1,2,2-Tetrachloroethane | 22.2 | | μg/l | | 20.0 | | 111 | 46-157 | 5 | 30 |
| Tetrachloroethene | 17.8 | | μg/l | | 20.0 | | 89 | 64-148 | 2 | 30 |
| 1 Stradition oction is | 20.0 | | μg/l | | 20.0 | | 100 | 70-130 | 2 | 30 |

| | | | | | Spike | Source | | %REC | | RPD |
|--|---|------|--------------|------|--------------|--------|------------|------------------|-----|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111291 - SW846 5030 Water | ·MS | | | | | | | | | |
| LCS Dup (8111291-BSD1) | | | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| 1,1,1-Trichloroethane | 19.3 | | μg/l | | 20.0 | | 97 | 52-162 | 3 | 30 |
| 1,1,2-Trichloroethane | 21.9 | | μg/l | | 20.0 | | 109 | 52-150 | 4 | 30 |
| Trichloroethene | 19.8 | | μg/l | | 20.0 | | 99 | 71-157 | 2 | 30 |
| Trichlorofluoromethane (Freon 11) | 19.2 | | μg/l | | 20.0 | | 96 | 17-181 | 2 | 30 |
| Vinyl chloride | 22.6 | | μg/l | | 20.0 | | 113 | 1-251 | 3 | 30 |
| m,p-Xylene | 42.8 | | μg/l | | 40.0 | | 107 | 70-130 | 3 | 30 |
| o-Xylene | 22.0 | | μg/l | | 20.0 | | 110 | 70-130 | 0.7 | 30 |
| Surrogate: 4-Bromofluorobenzene | 32.0 | | μg/l | | 30.0 | | 106 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.0 | | μg/l | | 30.0 | | 100 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 Surrogate: Dibromofluoromethane | 31.1 32.0 | | μg/l μg/l | | 30.0 30.0 | | 104 107 | 70-130 70-130 | | |
| Matrix Spike (8111291-MS1) | Source: SA8750 | 1-01 | P3 | | 00.0 | | | | | |
| Prepared & Analyzed: 18-Nov-08 | 2 | - • | | | | | | | | |
| Benzene | 23.0 | | μg/l | | 20.0 | BRL | 115 | 70-130 | | |
| Bromodichloromethane | 29.2 | | μg/l | | 20.0 | BRL | 146 | 35-155 | | |
| Bromoform | 22.7 | | μg/l | | 20.0 | BRL | 113 | 45-169 | | |
| Bromomethane | 24.8 | | μg/l | | 20.0 | BRL | 124 | 1-242 | | |
| Carbon tetrachloride | 28.8 | QM7 | μg/l | | 20.0 | BRL | 144 | 70-140 | | |
| Chlorobenzene | 23.8 | | μg/l | | 20.0 | BRL | 119 | 70-130 | | |
| Chloroethane | 25.0 | | μg/l | | 20.0 | BRL | 125 | 14-230 | | |
| Chloroform | 27.4 | | μg/l | | 20.0 | BRL | 137 | 51-138 | | |
| Chloromethane | 20.8 | | μg/l | | 20.0 | BRL | 104 | 1-273 | | |
| Dibromochloromethane | 23.8 | | μg/l | | 20.0 | BRL | 119 | 53-149 | | |
| 1,2-Dichlorobenzene | 21.4 | | μg/l | | 20.0 | BRL | 107 | 18-190 | | |
| 1,3-Dichlorobenzene | 28.0 | | μg/l | | 20.0 | BRL | 140 | 59-156 | | |
| 1,4-Dichlorobenzene | 21.9 | | μg/l | | 20.0 | BRL | 109 | 18-190 | | |
| 1,1-Dichloroethane | 24.1 | | μg/l | | 20.0 | BRL | 121 | 59-155 | | |
| 1,2-Dichloroethane | 25.8 | | μg/l | | 20.0 | BRL | 129 | 49-155 | | |
| 1,1-Dichloroethene | 21.9 | | μg/l | | 20.0 | BRL | 110 | 70-130 | | |
| trans-1,2-Dichloroethene | 23.8 | | μg/l | | 20.0 | BRL | 119 | 54-156 | | |
| 1,2-Dichloropropane | 23.4 | | μg/l | | 20.0 | BRL | 117 | 1-210 | | |
| cis-1,3-Dichloropropene | 19.3 | | μg/l | | 20.0 | BRL | 97 | 1-227 | | |
| trans-1,3-Dichloropropene | 20.6 | | μg/l | | 20.0 | BRL | 103 | 17-183 | | |
| Ethylbenzene | 23.8 | | μg/l | | 20.0 | BRL | 119 | 37-162 | | |
| Methylene chloride | 25.8 | | μg/l | | 20.0 | BRL | 129 | 1-221 | | |
| 1,1,2,2-Tetrachloroethane | 30.4 | | μg/l | | 20.0 | BRL | 152 | 46-157 | | |
| Tetrachloroethene | 22.8 | | μg/l | | 20.0 | BRL | 114 | 64-148 | | |
| Toluene | 24.3 | | μg/l | | 20.0 | BRL | 122 | 70-130 | | |
| 1,1,1-Trichloroethane | 24.2 | | μg/l | | 20.0 | BRL | 121 | 52-162 | | |
| 1,1,2-Trichloroethane | 26.7 | | μg/l | | 20.0 | BRL | 134 | 52-150 | | |
| Trichloroethene | 23.9 | | μg/l | | 20.0 | BRL | 120 | 71-157 | | |
| Trichlorofluoromethane (Freon 11) | 24.6 | | μg/l | | 20.0 | BRL | 123 | 17-181 | | |
| Vinyl chloride | 19.4 | | μg/l | | 20.0 | BRL | 97 | 1-251 | | |
| Surrogate: 4-Bromofluorobenzene | 34.6 | | μg/l | | 30.0 | | 115 | 70-130 | | |
| Surrogate: Toluene-d8 | 31.1 | | μg/l | | 30.0 | | 104 105 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 Surrogate: Dibromofluoromethane | 37.6 35.0 | | µg/l µg/l | | 30.0 30.0 | | 125 117 | 70-130 70-130 | | |
| Matrix Spike Dup (8111291-MSD1) | Source: SA8750 | 1-01 | . 2 | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| Benzene | 26.5 | QC1 | μg/l | | 20.0 | BRL | 132 | 70-130 | 14 | 30 |

| | | | | | Spike | Source | | %REC | | RPD |
|-----------------------------------|----------------|------|-------|------|-------|--------|------|--------|-----|------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limi |
| Batch 8111291 - SW846 5030 Water | · MS | | | | | | | | | |
| Matrix Spike Dup (8111291-MSD1) | Source: SA8750 | 1-01 | | | | | | | | |
| Prepared & Analyzed: 18-Nov-08 | | | | | | | | | | |
| Bromodichloromethane | 40.1 | QC1 | μg/l | | 20.0 | BRL | 201 | 35-155 | 32 | 30 |
| Bromoform | 35.2 | QC1 | μg/l | | 20.0 | BRL | 176 | 45-169 | 43 | 30 |
| Bromomethane | 24.8 | | μg/l | | 20.0 | BRL | 124 | 1-242 | 0.2 | 30 |
| Carbon tetrachloride | 31.1 | QM7 | μg/l | | 20.0 | BRL | 156 | 70-140 | 8 | 30 |
| Chlorobenzene | 34.5 | QC1 | μg/l | | 20.0 | BRL | 172 | 70-130 | 37 | 30 |
| Chloroethane | 24.4 | | μg/l | | 20.0 | BRL | 122 | 14-230 | 3 | 30 |
| Chloroform | 31.7 | QC1 | μg/l | | 20.0 | BRL | 158 | 51-138 | 15 | 30 |
| Chloromethane | 20.6 | | μg/l | | 20.0 | BRL | 103 | 1-273 | 0.9 | 30 |
| Dibromochloromethane | 35.7 | QC1 | μg/l | | 20.0 | BRL | 179 | 53-149 | 40 | 30 |
| 1,2-Dichlorobenzene | 36.5 | QC1 | μg/l | | 20.0 | BRL | 182 | 18-190 | 52 | 30 |
| 1,3-Dichlorobenzene | 41.8 | QC1 | μg/l | | 20.0 | BRL | 209 | 59-156 | 40 | 30 |
| 1,4-Dichlorobenzene | 35.2 | QC1 | μg/l | | 20.0 | BRL | 176 | 18-190 | 47 | 30 |
| 1,1-Dichloroethane | 25.4 | | μg/l | | 20.0 | BRL | 127 | 59-155 | 5 | 30 |
| 1,2-Dichloroethane | 30.3 | | μg/l | | 20.0 | BRL | 152 | 49-155 | 16 | 30 |
| 1,1-Dichloroethene | 21.8 | | μg/l | | 20.0 | BRL | 109 | 70-130 | 0.7 | 30 |
| trans-1,2-Dichloroethene | 23.9 | | μg/l | | 20.0 | BRL | 120 | 54-156 | 0.6 | 30 |
| 1,2-Dichloropropane | 30.2 | | μg/l | | 20.0 | BRL | 151 | 1-210 | 25 | 30 |
| cis-1,3-Dichloropropene | 28.3 | QC1 | μg/l | | 20.0 | BRL | 141 | 1-227 | 38 | 30 |
| trans-1,3-Dichloropropene | 32.1 | QC1 | μg/l | | 20.0 | BRL | 161 | 17-183 | 43 | 30 |
| Ethylbenzene | 37.0 | QC1 | μg/l | | 20.0 | BRL | 185 | 37-162 | 43 | 30 |
| Methylene chloride | 26.3 | | μg/l | | 20.0 | BRL | 131 | 1-221 | 2 | 30 |
| 1,1,2,2-Tetrachloroethane | 45.6 | QC1 | μg/l | | 20.0 | BRL | 228 | 46-157 | 40 | 30 |
| Tetrachloroethene | 29.1 | | μg/l | | 20.0 | BRL | 146 | 64-148 | 24 | 30 |
| Toluene | 31.4 | QC1 | μg/l | | 20.0 | BRL | 157 | 70-130 | 25 | 30 |
| 1,1,1-Trichloroethane | 28.8 | | μg/l | | 20.0 | BRL | 144 | 52-162 | 18 | 30 |
| 1,1,2-Trichloroethane | 38.9 | QC1 | μg/l | | 20.0 | BRL | 195 | 52-150 | 37 | 30 |
| Trichloroethene | 28.8 | | μg/l | | 20.0 | BRL | 144 | 71-157 | 18 | 30 |
| Trichlorofluoromethane (Freon 11) | 23.0 | | μg/l | | 20.0 | BRL | 115 | 17-181 | 7 | 30 |
| Vinyl chloride | 17.8 | | μg/l | | 20.0 | BRL | 89 | 1-251 | 9 | 30 |
| Surrogate: 4-Bromofluorobenzene | 33.2 | | μg/l | | 30.0 | DILL | 111 | 70-130 | | |
| Surrogate: Toluene-d8 | 30.7 | | μg/l | | 30.0 | | 102 | 70-130 | | |
| Surrogate: 1,2-Dichloroethane-d4 | 33.3 | | μg/l | | 30.0 | | 111 | 70-130 | | |
| Surrogate: Dibromofluoromethane | 32.0 | | μg/l | | 30.0 | | 107 | 70-130 | | |

Soluble Metals by EPA 6000/7000 Series Methods - Quality Control

| | | | | Spike | Source | | %REC | | RPD |
|--|--------------------|-----------|--------------|--------------|--------|------------|------------------|-----|-------|
| Analyte(s) | Result F | lag Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111042 - SW846 3005A | | | | | | | | | |
| Blank (8111042-BLK1) | | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | |
| Thallium | BRL | μg/l | 5.00 | | | | | | |
| Zinc | BRL | μg/l | 7.50 | | | | | | |
| Antimony | BRL | μg/l | 6.00 | | | | | | |
| Lead | BRL | μg/l | 7.50 | | | | | | |
| Nickel | BRL | μg/l | 5.00 | | | | | | |
| Selenium | BRL | μg/l | 15.0 | | | | | | |
| Beryllium | BRL | μg/l | 2.00 | | | | | | |
| Cadmium | BRL | μg/l | 2.50 | | | | | | |
| Chromium | BRL | μg/l | 5.00 | | | | | | |
| Arsenic | BRL | μg/l | 4.00 | | | | | | |
| Vanadium | BRL | μg/l | 5.00 | | | | | | |
| Silver | BRL | μg/l | 5.00 | | | | | | |
| Barium | BRL | μg/l | 5.00 | | | | | | |
| I CS (9444042 BS4) | | | | | | | | | |
| LCS (8111042-BS1) Prepared & Analyzed: 21-Nov-08 | | | | | | | | | |
| Zinc | 1320 | | 7.50 | 1050 | | 105 | OE 11E | | |
| Nickel | | μg/l | | 1250 | | 105 | 85-115 85-115 | | |
| | 1250 | μg/l | 5.00 | 1250 | | 100 | 85-115 85-115 | | |
| Lead | 1260 | μg/l | 7.50 | 1250 | | 100 | 85-115 85-115 | | |
| Thallium | 1210 1230 | μg/l | 5.00 | 1250 | | 97 99 | 85-115 85-115 | | |
| Selenium | | μg/l | 15.0 | 1250 | | | 85-115 85-115 | | |
| Antimony Barium | 1310 1230 | μg/l | 6.00 5.00 | 1250 1250 | | 105 99 | 85-115 85-115 | | |
| | | μg/l | | | | | | | |
| Cadmium Chromium | 1330 1250 | μg/l | 2.50 5.00 | 1250 1250 | | 106 100 | 85-115 85-115 | | |
| | | μg/l | | | | | | | |
| Arsenic Beryllium | 1240 | μg/l | 4.00 | 1250 | | 99 | 85-115 85-115 | | |
| Vanadium | 1230 1220 | μg/l | 2.00 | 1250 1250 | | 99 98 | 85-115 85-115 | | |
| Silver | 1250 | μg/l | 5.00 5.00 | 1250 | | 100 | 85-115 85-115 | | |
| Silver | 1250 | μg/l | 5.00 | 1250 | | 100 | 00-110 | | |
| LCS Dup (8111042-BSD1) | | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | |
| Lead | 1200 | μg/l | 7.50 | 1250 | | 96 | 85-115 | 4 | 20 |
| Thallium | 1170 | μg/l | 5.00 | 1250 | | 94 | 85-115 | 3 | 20 |
| Antimony | 1270 | μg/l | 6.00 | 1250 | | 102 | 85-115 | 3 | 20 |
| Selenium | 1200 | μg/l | 15.0 | 1250 | | 96 | 85-115 | 3 | 20 |
| Nickel | 1220 | μg/l | 5.00 | 1250 | | 97 | 85-115 | 3 | 20 |
| Zinc | 1280 | μg/l | 7.50 | 1250 | | 102 | 85-115 | 3 | 20 |
| Barium | 1200 | μg/l | 5.00 | 1250 | | 96 | 85-115 | 3 | 20 |
| Silver | 1210 | μg/l | 5.00 | 1250 | | 97 | 85-115 | 3 | 20 |
| Chromium | 1210 | μg/l | 5.00 | 1250 | | 97 | 85-115 | 3 | 20 |
| Vanadium | 1180 | μg/l | 5.00 | 1250 | | 95 | 85-115 | 4 | 20 |
| Beryllium | 1190 | μg/l | 2.00 | 1250 | | 95 | 85-115 | 3 | 20 |
| Cadmium | 1290 | μg/l | 2.50 | 1250 | | 103 | 85-115 | 3 | 20 |
| Arsenic | 1200 | μg/l | 4.00 | 1250 | | 96 | 85-115 | 3 | 20 |
| Duplicate (8111042-DUP1) | Source: SA87401-01 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | |
| Thallium | 3.80 | J μg/l | 5.00 | | BRL | | | | 20 |
| Zinc | 22.6 | μg/l | 7.50 | | 23.7 | | | 5 | 20 |
| Selenium | BRL | μg/l | 15.0 | | BRL | | | | 20 |

Soluble Metals by EPA 6000/7000 Series Methods - Quality Control

| | | | | | Spike | Source | | %REC | | RPD |
|---------------------------------|----------------|------|-------|------|-------|--------|------|--------|------|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111042 - SW846 3005A | | | | | | | | | | |
| <u>Duplicate (8111042-DUP1)</u> | Source: SA8740 | 1-01 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | | |
| Antimony | BRL | | μg/l | 6.00 | | BRL | | | | 20 |
| Nickel | BRL | | μg/l | 5.00 | | BRL | | | | 20 |
| Lead | 12.6 | QR8 | μg/l | 7.50 | | 6.00 | | | 71 | 20 |
| Barium | 8.20 | | μg/l | 5.00 | | 7.90 | | | 3 | 20 |
| Silver | BRL | | μg/l | 5.00 | | BRL | | | | 20 |
| Beryllium | BRL | | μg/l | 2.00 | | BRL | | | | 20 |
| Chromium | BRL | | μg/l | 5.00 | | BRL | | | | 20 |
| Arsenic | BRL | | μg/l | 4.00 | | BRL | | | | 20 |
| Cadmium | 0.500 | J | μg/l | 2.50 | | BRL | | | | 20 |
| Vanadium | BRL | | μg/l | 5.00 | | BRL | | | | 20 |
| Matrix Spike (8111042-MS1) | Source: SA8740 | 1-02 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | | |
| Antimony | 1300 | | μg/l | 6.00 | 1250 | BRL | 104 | 75-125 | | |
| Zinc | 1310 | | μg/l | 7.50 | 1250 | 16.9 | 103 | 75-125 | | |
| Thallium | 1190 | | μg/l | 5.00 | 1250 | BRL | 95 | 75-125 | | |
| Selenium | 1220 | | μg/l | 15.0 | 1250 | BRL | 97 | 75-125 | | |
| Lead | 1230 | | μg/l | 7.50 | 1250 | 5.35 | 98 | 75-125 | | |
| Nickel | 1230 | | μg/l | 5.00 | 1250 | 4.50 | 98 | 75-125 | | |
| Beryllium | 1210 | | μg/l | 2.00 | 1250 | BRL | 97 | 75-125 | | |
| Chromium | 1220 | | μg/l | 5.00 | 1250 | BRL | 97 | 75-125 | | |
| Vanadium | 1200 | | μg/l | 5.00 | 1250 | BRL | 96 | 70-130 | | |
| Silver | 1230 | | μg/l | 5.00 | 1250 | BRL | 98 | 75-125 | | |
| Arsenic | 1210 | | μg/l | 4.00 | 1250 | BRL | 97 | 75-125 | | |
| Barium | 1260 | | μg/l | 5.00 | 1250 | 69.4 | 95 | 75-125 | | |
| Cadmium | 1300 | | μg/l | 2.50 | 1250 | BRL | 104 | 75-125 | | |
| Matrix Spike Dup (8111042-MSD1) | Source: SA8740 | 1-02 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | | |
| Nickel | 1230 | | μg/l | 5.00 | 1250 | 4.50 | 98 | 75-125 | 0.3 | 20 |
| Zinc | 1310 | | μg/l | 7.50 | 1250 | 16.9 | 103 | 75-125 | 0.3 | 20 |
| Thallium | 1180 | | μg/l | 5.00 | 1250 | BRL | 95 | 75-125 | 0.5 | 20 |
| Selenium | 1220 | | μg/l | 15.0 | 1250 | BRL | 97 | 75-125 | 0.04 | 20 |
| Lead | 1220 | | μg/l | 7.50 | 1250 | 5.35 | 97 | 75-125 | 0.6 | 20 |
| Antimony | 1290 | | μg/l | 6.00 | 1250 | BRL | 103 | 75-125 | 0.4 | 20 |
| Arsenic | 1210 | | μg/l | 4.00 | 1250 | BRL | 97 | 75-125 | 0 | 20 |
| Silver | 1230 | | μg/l | 5.00 | 1250 | BRL | 98 | 75-125 | 0 | 20 |
| Cadmium | 1300 | | μg/l | 2.50 | 1250 | BRL | 104 | 75-125 | 0.5 | 20 |
| Barium | 1260 | | μg/l | 5.00 | 1250 | 69.4 | 95 | 75-125 | 0.04 | 20 |
| Chromium | 1220 | | μg/l | 5.00 | 1250 | BRL | 98 | 75-125 | 0.6 | 20 |
| Vanadium | 1200 | | μg/l | 5.00 | 1250 | BRL | 96 | 70-130 | 0.5 | 20 |
| Beryllium | 1200 | | μg/l | 2.00 | 1250 | BRL | 96 | 75-125 | 0.8 | 20 |
| Post Spike (8111042-PS1) | Source: SA8740 | 1-02 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | | |
| Thallium | 1250 | | μg/l | 5.00 | 1250 | BRL | 100 | 80-120 | | |
| Nickel | 1270 | | μg/l | 5.00 | 1250 | 4.50 | 101 | 80-120 | | |
| Lead | 1280 | | μg/l | 7.50 | 1250 | 5.35 | 102 | 80-120 | | |
| Zinc | 1350 | | μg/l | 7.50 | 1250 | 16.9 | 107 | 80-120 | | |
| Antimony | 1320 | | μg/l | 6.00 | 1250 | BRL | 106 | 80-120 | | |
| Selenium | 1260 | | μg/l | 15.0 | 1250 | BRL | 101 | 80-120 | | |

Soluble Metals by EPA 6000/7000 Series Methods - Quality Control

| | | | | | Spike | Source | | %REC | | RPD |
|--|--|-----------|--------------|-------------|------------|------------|-----------|------------------|-----|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Batch 8111042 - SW846 3005A | | | | | | | | | | |
| Post Spike (8111042-PS1) | Source: SA87401 | 1-02 | | | | | | | | |
| Prepared & Analyzed: 21-Nov-08 | | | | | | | | | | |
| Barium | 1320 | | μg/l | 5.00 | 1250 | 69.4 | 100 | 80-120 | | |
| Chromium | 1260 | | μg/l | 5.00 | 1250 | BRL | 101 | 80-120 | | |
| Cadmium | 1350 | | μg/l | 2.50 | 1250 | BRL | 108 | 80-120 | | |
| Arsenic | 1250 | | μg/l | 4.00 | 1250 | BRL | 100 | 80-120 | | |
| Beryllium | 1260 | | μg/l | 2.00 | 1250 | BRL | 100 | 80-120 | | |
| Vanadium | 1240 | | μg/l | 5.00 | 1250 | BRL | 99 | 80-120 | | |
| Silver | 1270 | | μg/l | 5.00 | 1250 | BRL | 102 | 80-120 | | |
| | Soluble Me | tals by l | EPA 200 S | Series Meth | ods - Qual | ity Contro | ol | | | |
| | | | | | Spike | Source | | %REC | | RPD |
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| Blank (8111043-BLK1) | . 00 | | | | | | | | | |
| <u> </u> | | | | | | | | | | |
| Prepared: 21-Nov-08 Analyzed: 24-Nov | v-us BRL | | | 0.000 | | | | | | |
| Mercury | DKL | | μg/l | 0.200 | | | | | | |
| LCS (8111043-BS1) | | | | | | | | | | |
| Prepared: 21-Nov-08 Analyzed: 24-Nov | /-08 | | | | | | | | | |
| Mercury | 4.48 | | μg/l | 0.200 | 5.00 | | 90 | 85-115 | | |
| Duplicate (8111043-DUP1) | Source: SA87371 | I-01 | | | | | | | | |
| | | | | | | | | | | |
| Prepared: 21-Nov-08 Analyzed: 24-Nov | <i>ı</i> -08 | | | | | | | | | |
| Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury | v-08 BRL | | μg/l | 0.200 | | BRL | | | | 20 |
| Mercury | BRL | 1-02 | µg/l | 0.200 | | BRL | | | | 20 |
| Mercury Matrix Spike (8111043-MS1) | BRL Source: SA87371 | 1-02 | μg/l | 0.200 | | BRL | | | | 20 |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov | BRL Source: SA87371 v-08 | 1-02 | | | 5.00 | | 100 | 75-125 | | 20 |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury | BRL Source: SA87374 v-08 4.99 | | µg/l µg/l | 0.200 | 5.00 | BRL BRL | 100 | 75-125 | | 20 |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury Matrix Spike Dup (8111043-MSD1) | BRL Source: SA87374 v-08 4.99 Source: SA87374 | | | | 5.00 | | 100 | 75-125 | | 20 |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury Matrix Spike Dup (8111043-MSD1) Prepared: 21-Nov-08 Analyzed: 24-Nov | BRL Source: SA87374 v-08 4.99 Source: SA87374 v-08 | | μg/l | 0.200 | | BRL | | | | |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury Matrix Spike Dup (8111043-MSD1) | BRL Source: SA87374 v-08 4.99 Source: SA87374 | | | | 5.00 | | 100 92 | 75-125 75-125 | 8 | 20 |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury Matrix Spike Dup (8111043-MSD1) Prepared: 21-Nov-08 Analyzed: 24-Nov | BRL Source: SA87374 v-08 4.99 Source: SA87374 v-08 | 1-02 | μg/l | 0.200 | | BRL | | | 8 | |
| Mercury Matrix Spike (8111043-MS1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury Matrix Spike Dup (8111043-MSD1) Prepared: 21-Nov-08 Analyzed: 24-Nov Mercury | BRL Source: SA87374 v-08 4.99 Source: SA87374 v-08 4.61 Source: SA87374 | 1-02 | μg/l | 0.200 | | BRL | | | 8 | |

Notes and Definitions

QC1 Analyte out of acceptance range.

QM7 The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.

QR2 The RPD result exceeded the QC control limits; however, both percent recoveries were acceptable. Sample results for the QC batch were accepted based on percent recoveries and completeness of QC data.

QR8 Analyses are not controlled on RPD values from sample concentrations that are less than 5 times the reporting level. The batch is accepted based upon the difference between the sample and duplicate is less than or equal to the reporting limit.

BRL Below Reporting Limit - Analyte NOT DETECTED at or above the reporting limit

dry Sample results reported on a dry weight basis

NR Not Reported

RPD Relative Percent Difference

J Detected but below the Reporting Limit; therefore, result is an estimated concentration (CLP J-Flag).

A plus sign (+) in the Method Reference column indicates the method is not accredited by NELAC.

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

<u>Matrix Spike</u>: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

Validated by: Hanibal C. Tayeh, Ph.D. Rebecca Merz

MADEP MCP ANALYTICAL METHOD REPORT CERTIFICATION FORM

| Labor | ratory Name: S | Spectrum Analytical, I | nc Agawam, MA | Λ | Project #: 127000 | 58 | | |
|---------|------------------------------|---|------------------------|-----------------------|------------------------|-----------------------|--------------|------|
| Projec | ct Location: W | alpole Park South-Wa | alpole, MA | | MADEP RTN 1: | | | |
| | form provides of 371-01 thro | certifications for the fugh SA87371-09 | ollowing data set: | | | | | |
| Samp | le matrices: | Aqueous Ground | l Water | | | | | |
| | | □ 8260B | □ 8151A | □ 8330 | ☑ 6010B | ☑ 7470A/1A | | |
| | SW-846 ods Used | □ 8270C | □ 8081A | □ VPH | □ 6020 | □ 9014M ² | | |
| 1,10011 | ous escu | □ 8082 | □ 8021B | □ ЕРН | □ 7000S ³ | □ 7196A | | |
| 2 M - S | SW-846 Method 9 | Number (RTN), if known 014 or MADEP Physiologie 000 Series List individual | | (PAC) Method | | | | |
| | | An affirmative respo | nse to questions A | , B, C and D is requ | ired for "Presumpt | ive Certainty" status | S | |
| A | | mples received by the astody documentation | | ndition consistent wi | th that described on | the | ☑ Yes | □ No |
| В | followed, in | A/QC procedures required actuding the requirement of the performance standard | ent to note and disc | | | | ☑ Yes | □ No |
| C | Certainty", | ta included in this rep as described in Sectio ssurance and Quality O Data"? | n 2.0 (a), (b), (c) an | nd (d) of the MADE | P document CAM V | VII A, | ☑ Yes | □ No |
| D | | ns (see Section 11.3 o | | | without significant | | ☐ Yes | □ No |
| | | A response to | questions E and F | below is required fo | or "Presumptive Ce | rtainty" status | | |
| E | Were all an achieved? | alytical QC performa | nce standards and r | ecommendations for | the specified method | ods | ☑ Yes | □ No |
| F | Were result | s for all analyte-list co | ompounds/elements | s for the specified m | ethod(s) reported? | | ☑ Yes | □ No |
| • | | All negative res | ponses are address | sed in a case narrati | ve on the cover pag | e of this report. | | |
| respo | nsible for obta | attest under the pair aining the informatio ief, accurate and com | n, the material co | | ytical report is, to t | | | |
| | | | | | Γ | Date: 11/26/2008 | | |



CHAIN OF CUSTODY RECORD

Page 1 of 1

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| Condition upon receipt: Elect Ambient EleC | EDD Format PDF, Excel wall Dis + in 49 | E-mail to raymond johnson & throtech com | Fax results when available to (| | TripBlank IIIIOS | GHC-6-GW | MW. 3- GW | MD-6-213 Of | MW-2-GW | MW.9-GW | R12-3- GW | N5-8-218 | R12-10-GW | Sample Id: | G=Grab C= | GW= Water X2=_ | 2=HCl 3=H ₂ SO ₄ 4=HNO ₃ 8= NaHSO ₄ 9=_ | Ray Johnson | MA | Grant | Took Distri | |
| | fall Dis + | of Hotel | | 3.5 | | | | | | | 30/11/11 | 4 | 11/11/08 | Date: | C=Composite | | 4=HNO ₃ 5=NaOH | | IOLIO | | | |
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| | | | | | |) | | | | | | | | | | | | 7 | - | 70 | 7 0 | NA Q |
| | 11/12/08 | 11/12/08 | Date: | | | | | | | | | かしん かまれる | * Medals have | State specific reporting standards: | Standard | □ Provide MA DEP MCP CAM Report □ Provide CTDEP RCP Report QA/QC Reporting Level | QA Reporting Notes: (check if needed) | towar | Sta | rek South | | |
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ANALYTICAL REPORT

Lab Number: L0818397

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: lan Cannan

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 Report Date: 12/23/08

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Project Number: 12700058 Report Date: 12/23/08

Sample Location

WALPOLE, MA

Alpha Sample ID

L0818397-01

Client ID

RIZ-10

Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Project Number: 12700058 Report Date: 12/23/08

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | NO |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | NO |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L0818397

Project Name: WALPOLE PARK SOUTH Lab Number:

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

MCP Related Narratives

Sample Receipt

In reference to question C:

The samples were received at the laboratory requiring filtration for Dissolved Metals; however, the samples were received beyond the 24 hour holding time recommended for filtration. The samples were filtered and preserved appropriately.

Dissolved Metals

L0818397-01 has an elevated detection limit for Thallium due to the dilution required by the high concentrations of non-target analytes. The requested reporting limit was achieved.



12230812:22

Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Case Narrative (continued)

In reference to question F:

All samples were analyzed for a subset of MCP elements per the Chain of Custody.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative

Unibeth & Simin

ALPHA

Date: 12/23/08

METALS



Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Project Number: 12700058 Report Date: 12/23/08

SAMPLE RESULTS

Lab ID: L0818397-01 Date Collected: 12/11/08 09:30

Client ID: RIZ-10 Date Received: 12/15/08
Sample Location: WALPOLE, MA Field Prep: Not Specified

Matrix: Water

Dilution Date Date Prep Analytical Method **Factor** Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst

Dissolved Metals by MCP 6000/7000 series

Thallium, Dissolved ND mg/l 0.0020 4 12/16/08 11:15 12/17/08 00:50 EPA 3005A 64,6020A BM



Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Project Number: 12700058 Report Date: 12/23/08

Method Blank Analysis Batch Quality Control

Dilution Date Date Analytical Method Analyst **Parameter Result Qualifier Units RDL Factor Prepared Analyzed** Dissolved Metals by MCP 6000/7000 series for sample(s): 01 Batch: WG347338-1 Thallium, Dissolved ND mg/l 0.0005 12/16/08 23:33 64,6020A ВМ 12/16/08 11:15

Prep Information

Digestion Method: EPA 3005A



Lab Control Sample Analysis Batch Quality Control

Lab Number:

L0818397

Project Number: 12700058

Project Name:

WALPOLE PARK SOUTH

Report Date:

12/23/08

| Parameter | LCS %Recovery | | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|--|-----------------------|----|-------------------|---------------------|-----|------------|
| Dissolved Metals by MCP 6000/7000 series | Associated sample(s): | 01 | Batch: WG347338 | -2 WG347338-3 | | |
| Thallium, Dissolved | 95 | | 96 | 80-120 | 1 | 20 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0818397

Project Number: 12700058 Report Date: 12/23/08

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal A Absent

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|----------------------------------|--------|----|------|------|--------|-------------------|
| L0818397-01A | Plastic 500ml unpreserved | А | 7 | 2C | Υ | Absent | - |
| L0818397-01B | Plastic 500ml HNO3 preserved spl | Α | <2 | 2C | Υ | Absent | MCP-TL-6020S(180) |



Project Name:WALPOLE PARK SOUTHLab Number:L0818397Project Number:12700058Report Date:12/23/08

GLOSSARY

Acronyms

- EPA Environmental Protection Agency.
- LCS Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.
- LCSD- Laboratory Control Sample Duplicate: Refer to LCS.
- MS Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.
- MSD Matrix Spike Sample Duplicate: Refer to MS.
- NA Not Applicable.
- NI Not Ignitable.
- NC Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.
- ND Not detected at the reported detection limit for the sample.
- RDL Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
- RPD Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

The following data qualifiers have been identified for use under the CT DEP Reasonable Confidence Protocols.

- A Spectra identified as "Aldol Condensation Product".
- B The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- J Estimated value. The analyte was tentatively identified; the quantitation is an estimation. (Tentatively identified compounds only.)

Standard Qualifiers

H - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.

Report Format: Data Usability Report



Project Name: WALPOLE PARK SOUTH Lab Number: L0818397
Project Number: 12700058 Report Date: 12/23/08

REFERENCES

Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



12230812:22 PLEASE ANSWER QUESTIONS ABOVE MA MCP or CT RCP? IS YOUR PROJECT ALPHA Lab ID DLs <RCGW-1 Other Project Specific Requirements/Comments/Detection Limits: Email: ian.cannan@tetratech.com Phone: 508-903-2039 Fax: 508-903-2001 Framingham, MA Client Information Westborough, MA These samples have been Previously analyzed by Alpha Address: One Grant Street Client: Tetra Tech Rizzo TEL: 508-898-9220 FAX: 508-898-9193 (Lab Use Only) TEL: 508-822-9300 Mansfield, MA FAX: 508-822-3288 RIZ-10 CHAIN OF CUSTODY Sample ID Due Date: 17/77 (%Time: Project Location: Walpole, MA Standard
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 Standard Project Manager: Ray Johnson Turn-Around Time Project #: 12700058 Project Name: Walpole Park South Project Information ALPHA Quote #: 12-11-08 Date Collection 0930 Time □ Rush (ONLY IF PRE-APPROVED) Relinguisbed By: PAGE Sample Matrix GW Container Type 읶 Preservative SC Sampler's Initials ס 12-15-08 12-15-02 Date Rec'd in Lab: \boxtimes ☐ Yes ⊠ Yes Regulatory Requirements/Report Limits ANALYSIS MCP PRESUMPTIVE CERTAINTY-CT REASONABLE CONFIDENCE PROTOCOLS ☐ FAX Report Information Data Deliverables MA MCP CAM State/Fed Program ⊠ ADE> Dissolved Thallium Date/Time □ 8 S S 2-115/08 Add'l Deliverables **⊠** EMAIL Are CT RCP (Reasonable Confidence Protocols) Required? Are MCP Analytical Methods Required? Criteria Billing Information ALPHA Job #: 208/8397 Same as Client info 2/15/68 Date/Time Please print clearly, legibly and completely. Samples can not be logged in and timaround time alock will negative and timaround time alock will negative and the logged in an and timaround time alock will negative. All samples submitted are subject to Alpha's Payment Terms. Sample Specific Comments □ Done PO # 🛛 Lab to do ☐ Not Needec Lab to filter (Please specif) below) 🛛 Lab to do Preservation **Filtration** SAMPLE HANDLING SELTIOB



ANALYTICAL REPORT

Lab Number: L0907670

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: Ray Johnson

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Report Date: 06/17/09

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Number: 12700053 **Report Date:** 06/17/09

| Alpha Sample ID | Client ID | Sample Location | Collection Date/Time |
|--------------------|-----------|--------------------|-------------------------|
| L0907670-01 | RIZ-10 | WALPOLE, MA | 06/10/09 13:00 |
| L0907670-02 | RIZ-8 | WALPOLE, MA | 06/10/09 13:28 |
| L0907670-03 | RIZ-8S | WALPOLE, MA | 06/10/09 13:18 |
| L0907670-04 | MW-9 | WALPOLE, MA | 06/10/09 13:45 |
| L0907670-05 | GHC-6 | WALPOLE, MA | 06/10/09 14:05 |
| L0907670-06 | RIZ-3 | WALPOLE, MA | 06/10/09 14:12 |
| L0907670-07 | MW-2 | WALPOLE, MA | 06/10/09 14:26 |
| L0907670-08 | RIZ-9 | WALPOLE, MA | 06/10/09 14:41 |

Project Number: 12700053 Report Date: 06/17/09

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | YES |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L0907670

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700053 **Report Date:** 06/17/09

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

For additional information, please contact Client Services at 800-624-9220.

MCP Related Narratives

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

Metals

L0907670-01 through -08 have elevated detection limits for Antimony and Thallium due to the dilutions required by the high concentrations of non-target analytes. The requested reporting limits were achieved.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Unabeth & Simuro

Authorized Signature:

Title: Technical Director/Representative

Date: 06/17/09

ORGANICS



VOLATILES



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: Report Date: 12700053 06/17/09

SAMPLE RESULTS

Lab ID: Date Collected: L0907670-01 06/10/09 13:00

Client ID: RIZ-10

Date Received: 06/11/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 06/12/09 14:10 Analytical Date:

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborough | h Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 Report Date: 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-01 Date Collected: 06/10/09 13:00

Client ID: RIZ-10 Date Received: 06/11/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL Parameter Qualifier Units **Dilution Factor** Result Volatile Organics by GC/MS - Westborough Lab ND 1,3-Dichlorobenzene ug/l 0.50 1 ND 0.50 1 1,4-Dichlorobenzene ug/l Styrene ND 0.50 1 ug/l o-Xylene ND ug/l 0.50 1 ND 1,1-Dichloropropene ug/l 0.50 1 ND 0.50 2,2-Dichloropropane ug/l 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l ND 1 1,2,3-Trichloropropane ug/l 0.50 Bromochloromethane ND ug/l 0.50 1 ND 0.50 n-Butylbenzene ug/l 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 ND 1 1,2,3-Trichlorobenzene ug/l 0.50 ND 0.50 1 1,2,4-Trichlorobenzene ug/l 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND 0.50 1 ug/l Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND 0.50 1 ug/l ND p-Chlorotoluene ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 ND 1,2-Dibromoethane ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 ND 1,3-Dichloropropane ug/l 0.50 1

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 105 | | 80-120 | |
| 4-Bromofluorobenzene | 83 | | 80-120 | |

ND



1

0.50

ug/l

Methyl tert butyl ether

06/11/09

Date Received:

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-02 Date Collected: 06/10/09 13:28

Client ID: RIZ-8

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/12/09 14:47

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbord | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-02 Date Collected: 06/10/09 13:28

Client ID: RIZ-8 Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| ' | | | | • | |
|-------------------------------------|------------|-----------|-------|------|----------------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Facto |
| Volatile Organics by GC/MS - Westbo | orough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 107 | | 80-120 | |
| 4-Bromofluorobenzene | 83 | | 80-120 | |



06/11/09

See Narrative

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-03 Date Collected: 06/10/09 13:18

Client ID: RIZ-8S

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/12/09 15:24

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | gh Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | 0.71 | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 Report Date: 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-03 Date Collected: 06/10/09 13:18

Client ID: RIZ-8S Date Received: 06/11/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL Parameter Qualifier Units **Dilution Factor** Result Volatile Organics by GC/MS - Westborough Lab ND 1,3-Dichlorobenzene ug/l 0.50 1 ND 1 1,4-Dichlorobenzene ug/l 0.50 ND 0.50 1 Styrene ug/l o-Xylene ND ug/l 0.50 1 ND 1,1-Dichloropropene ug/l 0.50 1 ND 0.50 2,2-Dichloropropane ug/l 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l ND 1 1,2,3-Trichloropropane ug/l 0.50 Bromochloromethane ND ug/l 0.50 1 ND n-Butylbenzene ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 ND 1 1,2,3-Trichlorobenzene ug/l 0.50 ND 0.50 1,2,4-Trichlorobenzene ug/l 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 ND 1,3,5-Trimethylbenzene 0.50 1 ug/l Bromobenzene ND ug/l 0.50 1 ND 0.50 1 o-Chlorotoluene ug/l p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 ND 1,2-Dibromoethane ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 ND 1,3-Dichloropropane ug/l 0.50 1 ND Methyl tert butyl ether 0.50 ug/l 1

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 106 | | 80-120 | |
| 4-Bromofluorobenzene | 84 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: Report Date: 12700053 06/17/09

SAMPLE RESULTS

Date Collected: Lab ID: L0907670-04 06/10/09 13:45

Client ID: MW-9

Date Received: 06/11/09 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 06/12/09 16:01

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbore | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | 0.75 | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-04 Date Collected: 06/10/09 13:45

Client ID: MW-9 Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Campio Ecodationi Witzi CEE, in | , · | | | .а. тор. | oo mana | |
|--------------------------------------|----------|-----------|-------|----------|-----------------|--|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | |
| Volatile Organics by GC/MS - Westbor | ough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 | |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 | |
| Styrene | ND | | ug/l | 0.50 | 1 | |
| o-Xylene | ND | | ug/l | 0.50 | 1 | |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 | |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 | |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 | |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 | |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 | |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 | |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 | |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 | |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 | |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 | |
| Naphthalene | ND | | ug/l | 0.50 | 1 | |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 | |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 | |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 | |
| Bromobenzene | ND | | ug/l | 0.50 | 1 | |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 | |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 | |
| Dibromomethane | ND | | ug/l | 0.50 | 1 | |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 | |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 | |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 | |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 | |
| | | | | | | |

| Surrogate | % Recovery | Acceptance % Recovery Qualifier Criteria | | | | |
|------------------------|------------|---|--------|--|--|--|
| 1,2-Dichlorobenzene-d4 | 106 | | 80-120 | | | |
| 4-Bromofluorobenzene | 84 | | 80-120 | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: Report Date: 12700053 06/17/09

SAMPLE RESULTS

Lab ID: Date Collected: L0907670-05 06/10/09 14:05

Client ID: GHC-6

Date Received: 06/11/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 06/12/09 16:38

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbord | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-05 Date Collected: 06/10/09 14:05

Client ID: GHC-6 Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| , | | | | • | |
|-------------------------------------|-----------|-----------|-------|------|-----------------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
| Volatile Organics by GC/MS - Westbo | rough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Surrogate | % Recovery | Acceptance ecovery Qualifier Criteria | | | |
|------------------------|------------|--|--------|--|--|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 | | |
| 4-Bromofluorobenzene | 84 | | 80-120 | | |



06/11/09

Date Received:

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-06 Date Collected: 06/10/09 14:12

Client ID: RIZ-3

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/12/09 17:15

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbord | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-06 Date Collected: 06/10/09 14:12

Client ID: RIZ-3 Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Surrogate | % Recovery | Qualifier | | |
|------------------------|------------|-----------|--------|--|
| 1,2-Dichlorobenzene-d4 | 107 | | 80-120 | |
| 4-Bromofluorobenzene | 83 | | 80-120 | |



06/11/09

See Narrative

Date Received:

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-07 Date Collected: 06/10/09 14:26

Client ID: MW-2

Sample Location: WALPOLE, MA Field Prep:

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/12/09 17:52

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborough | n Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | 1.0 | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 Report Date: 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-07 Date Collected: 06/10/09 14:26

Client ID: MW-2 Date Received: 06/11/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL Parameter Qualifier Units **Dilution Factor** Result Volatile Organics by GC/MS - Westborough Lab 1,3-Dichlorobenzene ND ug/l 0.50 1 ND 1 1,4-Dichlorobenzene ug/l 0.50 ND 0.50 1 Styrene ug/l o-Xylene ND ug/l 0.50 1 ND 1,1-Dichloropropene ug/l 0.50 1 ND 0.50 2,2-Dichloropropane ug/l 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l ND 1,2,3-Trichloropropane ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 ND n-Butylbenzene ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 ND 1,2,3-Trichlorobenzene ug/l 0.50 1 ND 0.50 1,2,4-Trichlorobenzene ug/l 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 ND 1,3,5-Trimethylbenzene 0.50 1 ug/l Bromobenzene ND ug/l 0.50 1 ND 0.50 1 o-Chlorotoluene ug/l ug/l p-Chlorotoluene ND 0.50 1 Dibromomethane ND ug/l 0.50 1 ND 1,2-Dibromoethane ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 ND 1,3-Dichloropropane ug/l 0.50 1 ND Methyl tert butyl ether 0.50 ug/l 1

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | | |
|------------------------|------------|-----------|------------------------|--|--|
| 1,2-Dichlorobenzene-d4 | 108 | | 80-120 | | |
| 4-Bromofluorobenzene | 84 | | 80-120 | | |



06/11/09

Date Received:

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-08 Date Collected: 06/10/09 14:41

Client ID: RIZ-9

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/12/09 18:29

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborough | Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-08 Date Collected: 06/10/09 14:41

Client ID: RIZ-9 Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------------|-----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbor | rough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 105 | | 80-120 | |
| 4-Bromofluorobenzene | 84 | | 80-120 | |



Project Number: 12700053 **Report Date:** 06/17/09

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/12/09 08:00

Analyst: TT

| rameter | Result | Qualifier | Units | RDL | |
|---------------------------|----------------|------------------|-------|--------|------------|
| latile Organics by GC/MS | Westborough La | b for sample(s): | 01-08 | Batch: | WG366434-4 |
| Methylene chloride | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | |
| Chloroform | ND | | ug/l | 0.50 | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | |
| Dibromochloromethane | ND | | ug/l | 0.50 | |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | |
| Tetrachloroethene | ND | | ug/l | 0.50 | |
| Chlorobenzene | ND | | ug/l | 0.50 | |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | |
| Bromodichloromethane | ND | | ug/l | 0.50 | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| Bromoform | ND | | ug/l | 0.50 | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | |
| Benzene | ND | | ug/l | 0.50 | |
| Toluene | ND | | ug/l | 0.50 | |
| Ethylbenzene | ND | | ug/l | 0.50 | |
| p/m-Xylene | ND | | ug/l | 0.50 | |
| Chloromethane | ND | | ug/l | 0.50 | |
| Bromomethane | ND | | ug/l | 0.50 | |
| Vinyl chloride | ND | | ug/l | 0.50 | |
| Chloroethane | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| Trichloroethene | ND | | ug/l | 0.50 | |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | |



Project Number: 12700053 **Report Date:** 06/17/09

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/12/09 08:00

Analyst: TT

| arameter | Result Qualifier | Units | RDL | |
|-----------------------------|---------------------------------|---------|--------|------------|
| olatile Organics by GC/MS | - Westborough Lab for sample(s) | : 01-08 | Batch: | WG366434-4 |
| 1,4-Dichlorobenzene | ND | ug/l | 0.50 | |
| Styrene | ND | ug/l | 0.50 | |
| o-Xylene | ND | ug/l | 0.50 | |
| 1,1-Dichloropropene | ND | ug/l | 0.50 | |
| 2,2-Dichloropropane | ND | ug/l | 0.50 | |
| 1,1,1,2-Tetrachloroethane | ND | ug/l | 0.50 | |
| 1,2,3-Trichloropropane | ND | ug/l | 0.50 | |
| Bromochloromethane | ND | ug/l | 0.50 | |
| n-Butylbenzene | ND | ug/l | 0.50 | |
| Dichlorodifluoromethane | ND | ug/l | 0.50 | |
| Hexachlorobutadiene | ND | ug/l | 0.50 | |
| Isopropylbenzene | ND | ug/l | 0.50 | |
| p-Isopropyltoluene | ND | ug/l | 0.50 | |
| Naphthalene | ND | ug/l | 0.50 | |
| n-Propylbenzene | ND | ug/l | 0.50 | |
| sec-Butylbenzene | ND | ug/l | 0.50 | |
| tert-Butylbenzene | ND | ug/l | 0.50 | |
| 1,2,3-Trichlorobenzene | ND | ug/l | 0.50 | |
| 1,2,4-Trichlorobenzene | ND | ug/l | 0.50 | |
| 1,2,4-Trimethylbenzene | ND | ug/l | 0.50 | |
| 1,3,5-Trimethylbenzene | ND | ug/l | 0.50 | |
| Bromobenzene | ND | ug/l | 0.50 | |
| o-Chlorotoluene | ND | ug/l | 0.50 | |
| p-Chlorotoluene | ND | ug/l | 0.50 | |
| Dibromomethane | ND | ug/l | 0.50 | |
| 1,2-Dibromoethane | ND | ug/l | 0.50 | |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 0.50 | |
| 1,3-Dichloropropane | ND | ug/l | 0.50 | |
| Methyl tert butyl ether | ND | ug/l | 0.50 | |
| | | | | |



Project Number: 12700053 **Report Date:** 06/17/09

Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/12/09 08:00

Analyst: TT

Parameter Result Qualifier Units RDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-08 Batch: WG366434-4

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | | Acceptance | | | | |
|------------------------|-----------|------------|----------|--|--|--|
| Surrogate | %Recovery | Qualifier | Criteria | | | |
| 1,2-Dichlorobenzene-d4 | 106 | | 80-120 | | | |
| 4-Bromofluorobenzene | 87 | | 80-120 | | | |



Lab Control Sample Analysis Batch Quality Control

WALPOLE PARK SOUTH

Project Number: 12700053

Project Name:

Lab Number:

L0907670

Report Date:

06/17/09

| rameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---|----------------------------|-------------------|---------------------|-----|------------|
| platile Organics by GC/MS - Westborough | h Lab Associated sample(s) | : 01-08 Batch: | WG366434-3 | | |
| Methylene chloride | 101 | - | 70-130 | - | |
| 1,1-Dichloroethane | 96 | - | 70-130 | - | |
| Chloroform | 99 | - | 70-130 | - | |
| Carbon tetrachloride | 91 | - | 70-130 | - | |
| 1,2-Dichloropropane | 97 | - | 70-130 | - | |
| Dibromochloromethane | 91 | - | 70-130 | - | |
| 1,1,2-Trichloroethane | 97 | - | 70-130 | - | |
| Tetrachloroethene | 94 | - | 70-130 | - | |
| Chlorobenzene | 97 | - | 70-130 | - | |
| Trichlorofluoromethane | 100 | - | 70-130 | - | |
| 1,2-Dichloroethane | 93 | - | 70-130 | - | |
| 1,1,1-Trichloroethane | 91 | - | 70-130 | - | |
| Bromodichloromethane | 94 | - | 70-130 | - | |
| trans-1,3-Dichloropropene | 85 | - | 70-130 | - | |
| cis-1,3-Dichloropropene | 85 | - | 70-130 | - | |
| Bromoform | 92 | - | 70-130 | - | |
| 1,1,2,2-Tetrachloroethane | 106 | - | 70-130 | - | |
| Benzene | 98 | - | 70-130 | - | |
| Toluene | 91 | - | 70-130 | - | |
| Ethylbenzene | 96 | - | 70-130 | - | |
| p/m-Xylene | 98 | - | 70-130 | - | |

Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

Report Date: 06/17/09

| rameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|---|--------------------------|-------------------|---------------------|-----|------------|
| platile Organics by GC/MS - Westborough L | ab Associated sample(s): | 01-08 Batch: | WG366434-3 | | |
| Chloromethane | 112 | - | 70-130 | - | |
| Bromomethane | 114 | - | 70-130 | - | |
| Vinyl chloride | 104 | - | 70-130 | - | |
| Chloroethane | 104 | - | 70-130 | - | |
| 1,1-Dichloroethene | 98 | - | 70-130 | - | |
| trans-1,2-Dichloroethene | 95 | - | 70-130 | - | |
| cis-1,2-Dichloroethene | 92 | - | 70-130 | - | |
| Trichloroethene | 86 | - | 70-130 | - | |
| 1,2-Dichlorobenzene | 97 | - | 70-130 | - | |
| 1,3-Dichlorobenzene | 96 | - | 70-130 | - | |
| 1,4-Dichlorobenzene | 95 | - | 70-130 | - | |
| Styrene | 96 | - | 70-130 | - | |
| o-Xylene | 91 | - | 70-130 | - | |
| 1,1-Dichloropropene | 92 | - | 70-130 | - | |
| 2,2-Dichloropropane | 94 | - | 70-130 | - | |
| 1,1,1,2-Tetrachloroethane | 95 | - | 70-130 | - | |
| 1,2,3-Trichloropropane | 97 | - | 70-130 | - | |
| Bromochloromethane | 100 | - | 70-130 | - | |
| n-Butylbenzene | 94 | - | 70-130 | - | |
| Dichlorodifluoromethane | 105 | - | 70-130 | - | |
| Hexachlorobutadiene | 100 | - | 70-130 | - | |



L0907670

Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

tty Control Lab Number:

Report Date: 06/17/09

| arameter | LCS %Recovery | LCSD %Recovery | %Recovery Limits | RPD | RPD Limits |
|-------------------------------|------------------------------------|-------------------|---------------------|-----|------------|
| olatile Organics by GC/MS - V | Westborough Lab Associated sample(| s): 01-08 Batch: | WG366434-3 | | |
| Isopropylbenzene | 96 | - | 70-130 | - | |
| p-Isopropyltoluene | 93 | - | 70-130 | - | |
| Naphthalene | 75 | - | 70-130 | - | |
| n-Propylbenzene | 96 | - | 70-130 | - | |
| sec-Butylbenzene | 96 | - | 70-130 | - | |
| tert-Butylbenzene | 94 | - | 70-130 | - | |
| 1,2,3-Trichlorobenzene | 89 | - | 70-130 | - | |
| 1,2,4-Trichlorobenzene | 90 | - | 70-130 | - | |
| 1,2,4-Trimethylbenzene | 89 | - | 70-130 | - | |
| 1,3,5-Trimethylbenzene | 87 | - | 70-130 | - | |
| Bromobenzene | 102 | - | 70-130 | - | |
| o-Chlorotoluene | 98 | - | 70-130 | - | |
| p-Chlorotoluene | 93 | - | 70-130 | - | |
| Dibromomethane | 94 | - | 70-130 | - | |
| 1,2-Dibromoethane | 94 | - | 70-130 | - | |
| 1,2-Dibromo-3-chloropropane | 103 | - | 70-130 | - | |
| 1,3-Dichloropropane | 92 | - | 70-130 | - | |
| Methyl tert butyl ether | 88 | - | 70-130 | - | |

Lab Control Sample Analysis Batch Quality Control

WALPOLE PARK SOUTH

Lab Number: L0907670

Project Number: Report Date: 06/17/09 12700053

LCS **LCSD** %Recovery Limits %Recovery %Recovery

RPD **RPD Limits** Parameter

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 Batch: WG366434-3

| Surrogate | LCS %Recovery Qualifier | LCSD %Recovery Qualifier | Acceptance Criteria |
|------------------------|----------------------------|-----------------------------|------------------------|
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 |
| 4-Bromofluorobenzene | 97 | | 80-120 |



Project Name:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|--------------------------------------|------------------|--------------|---------------|-----------------|---------------|------------------|--------------------|--------|---------------|
| /olatile Organics by GC/MS Sample | - Westborough La | b Associated | sample(s): 01 | -08 QC Ba | ch ID: WG3664 | 34-5 QC Sar | mple: L0907641 | I-01 C | Client ID: MS |
| Methylene chloride | ND | 4 | 4.1 | 103 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| Chloroform | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Carbon tetrachloride | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| Dibromochloromethane | ND | 4 | 3.6 | 90 | - | - | 70-130 | - | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 3.8 | 95 | - | - | 70-130 | - | 20 |
| Tetrachloroethene | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Chlorobenzene | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |
| Trichlorofluoromethane | ND | 4 | 4.4 | 109 | - | - | 70-130 | - | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Bromodichloromethane | ND | 4 | 3.8 | 95 | - | - | 70-130 | - | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 3.0 | 75 | - | - | 70-130 | - | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 3.7 | 93 | - | - | 70-130 | - | 20 |
| Bromoform | ND | 4 | 3.5 | 88 | - | - | 70-130 | - | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 4.1 | 103 | - | - | 70-130 | - | 20 |
| Benzene | ND | 4 | 4.2 | 105 | - | - | 70-130 | - | 20 |
| Toluene | ND | 4 | 3.8 | 96 | - | - | 70-130 | - | 20 |
| Ethylbenzene | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| p/m-Xylene | ND | 8 | 7.9 | 99 | - | - | 70-130 | - | 20 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

| arameter | Native Sample | MS Added | MS Found | | S overy | MSD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|---------------------------------|------------------|---------------|---------------|-------|------------|---------------|------------------|--------------------|-------|---------------|
| olatile Organics by GC/MS ample | - Westborough La | ab Associated | sample(s): 01 | -08 C | QC Batcl | n ID: WG36643 | 34-5 QC Sar | mple: L0907641 | -01 C | Client ID: MS |
| Chloromethane | ND | 4 | 4.2 | | 104 | | - | 70-130 | - | 20 |
| Bromomethane | ND | 4 | 4.7 | | 117 | - | - | 70-130 | - | 20 |
| Vinyl chloride | ND | 4 | 5.2 | | 131 | - | - | 70-130 | - | 20 |
| Chloroethane | ND | 4 | 4.6 | | 115 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.4 | | 111 | - | - | 70-130 | - | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.2 | | 105 | - | - | 70-130 | - | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 3.9 | | 99 | - | - | 70-130 | - | 20 |
| Trichloroethene | ND | 4 | 3.8 | | 96 | - | - | 70-130 | - | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 3.8 | | 95 | - | - | 70-130 | - | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 3.8 | | 96 | - | - | 70-130 | - | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 3.7 | | 93 | - | - | 70-130 | - | 20 |
| Styrene | ND | 4 | 3.8 | | 94 | - | - | 70-130 | - | 20 |
| o-Xylene | ND | 4 | 3.7 | | 94 | - | - | 70-130 | - | 20 |
| 1,1-Dichloropropene | ND | 4 | 3.8 | | 96 | - | - | 70-130 | - | 20 |
| 2,2-Dichloropropane | ND | 4 | 4.1 | | 103 | - | - | 70-130 | - | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 3.7 | | 92 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 3.8 | | 94 | - | - | 70-130 | - | 20 |
| Bromochloromethane | ND | 4 | 4.0 | | 101 | - | - | 70-130 | - | 20 |
| n-Butylbenzene | ND | 4 | 3.9 | | 98 | - | - | 70-130 | - | 20 |
| Dichlorodifluoromethane | ND | 4 | 3.8 | | 96 | - | - | 70-130 | - | 20 |
| Hexachlorobutadiene | ND | 4 | 4.1 | | 102 | - | - | 70-130 | - | 20 |
| | | | | | | | | | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Found | MSD %Recovery | Recovery Limits | RPD | RPD Limits |
|--------------------------------------|------------------|---------------|----------------|-----------------|---------------|------------------|--------------------|--------|--------------|
| Volatile Organics by GC/MS Sample | - Westborough La | ab Associated | sample(s): 01- | 08 QC Bato | h ID: WG36643 | 4-5 QC Sar | nple: L0907641 | I-01 C | lient ID: MS |
| Isopropylbenzene | ND | 4 | 3.5 | 88 | - | - | 70-130 | - | 20 |
| p-Isopropyltoluene | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| Naphthalene | ND | 4 | 3.0 | 76 | - | - | 70-130 | - | 20 |
| n-Propylbenzene | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| sec-Butylbenzene | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| tert-Butylbenzene | ND | 4 | 3.9 | 97 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.4 | 86 | - | - | 70-130 | - | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.6 | 89 | - | - | 70-130 | - | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 3.6 | 91 | - | - | 70-130 | - | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 3.5 | 87 | - | - | 70-130 | - | 20 |
| Bromobenzene | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| o-Chlorotoluene | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| p-Chlorotoluene | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| Dibromomethane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| 1,2-Dibromoethane | ND | 4 | 3.7 | 92 | - | - | 70-130 | - | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 3.4 | 86 | - | - | 70-130 | - | 20 |
| 1,3-Dichloropropane | ND | 4 | 3.6 | 90 | - | - | 70-130 | - | 20 |
| Methyl tert butyl ether | ND | 4 | 3.5 | 88 | - | - | 70-130 | - | 20 |
| | | | | | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number:

L0907670

Report Date:

06/17/09

| MS | MSD | Recovery |
|----|-----|----------|
| | | |

Parameter Native Sample MS Added MS Found %Recovery MSD Found %Recovery Limits RPD RPD Limits

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 QC Batch ID: WG366434-5 QC Sample: L0907641-01 Client ID: MS Sample

| | MS | MSD | Acceptance | |
|------------------------|----------------------|----------------------|------------|--|
| Surrogate | % Recovery Qualifier | % Recovery Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 96 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number:

L0907670

Report Date:

06/17/09

| arameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------------|------------------------|-------------|--------------|-------------------|
| olatile Organics by GC/MS - Westborough Lab | Associated sample(s): 01-08 | QC Batch ID: WG366434- | 6 QC Sample | e: L0907641- | 02 Client ID: DUP |
| Methylene chloride | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethane | ND | ND | ug/l | NC | 20 |
| Chloroform | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene | ND | ND | ug/l | NC | 20 |
| Chlorobenzene | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| Bromoform | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| Benzene | ND | ND | ug/l | NC | 20 |
| Toluene | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number:

L0907670

Report Date:

06/17/09

| arameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------------|------------------------|--------------|----------------|----------------|
| olatile Organics by GC/MS - Westborough Lab | Associated sample(s): 01-08 | QC Batch ID: WG366434- | -6 QC Sample | e: L0907641-02 | Client ID: DUP |
| Ethylbenzene | ND | ND | ug/l | NC | 20 |
| p/m-Xylene | ND | ND | ug/l | NC | 20 |
| Chloromethane | ND | ND | ug/l | NC | 20 |
| Bromomethane | ND | ND | ug/l | NC | 20 |
| Vinyl chloride | ND | ND | ug/l | NC | 20 |
| Chloroethane | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethene | ND | ND | ug/l | NC | 20 |
| trans-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| cis-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| Trichloroethene | ND | ND | ug/l | NC | 20 |
| 1,2-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,3-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,4-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| Styrene | ND | ND | ug/l | NC | 20 |
| o-Xylene | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloropropene | ND | ND | ug/l | NC | 20 |
| 2,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| 1,1,1,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichloropropane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------------|----------------------|-------------|------------------|------------------|
| Volatile Organics by GC/MS - Westborough Lab Sample | Associated sample(s): 01-08 | QC Batch ID: WG36643 | 34-6 QC Sar | mple: L0907641-0 | 2 Client ID: DUP |
| Bromochloromethane | ND | ND | ug/l | NC | 20 |
| n-Butylbenzene | ND | ND | ug/l | NC | 20 |
| Dichlorodifluoromethane | ND | ND | ug/l | NC | 20 |
| Hexachlorobutadiene | ND | ND | ug/l | NC | 20 |
| Isopropylbenzene | ND | ND | ug/l | NC | 20 |
| p-Isopropyltoluene | ND | ND | ug/l | NC | 20 |
| Naphthalene | ND | ND | ug/l | NC | 20 |
| n-Propylbenzene | ND | ND | ug/l | NC | 20 |
| sec-Butylbenzene | ND | ND | ug/l | NC | 20 |
| tert-Butylbenzene | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| 1,3,5-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| Bromobenzene | ND | ND | ug/l | NC | 20 |
| o-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| p-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| Dibromomethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dibromoethane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number:

L0907670

Report Date:

06/17/09

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|----------------------------|-----------------------|-------------|----------------|-------------------|
| Volatile Organics by GC/MS - Westborough Lab As Sample | ssociated sample(s): 01-08 | QC Batch ID: WG366434 | 1-6 QC Samp | ole: L0907641- | 02 Client ID: DUP |
| 1,2-Dibromo-3-chloropropane | ND | ND | ug/l | NC | 20 |
| 1,3-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Methyl tert butyl ether | ND | ND | ug/l | NC | 20 |

| | | | | | Acceptance | |
|------------------------|-----------|-----------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 104 | | 106 | | 80-120 | |
| 4-Bromofluorobenzene | 85 | | 85 | | 80-120 | |



METALS



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-01 Date Collected: 06/10/09 13:00

Client ID: Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:05 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.148 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 10:55 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:05 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:18 | EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-02

Client ID: RIZ-8

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 06/10/09 13:28

Date Received: 06/11/09

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:34 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.026 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 10:56 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:34 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:24 | EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-03 Date Collected: 06/10/09 13:18

Client ID: RIZ-8S Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:40 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.051 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 10:58 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:40 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:26 | EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-04

Client ID: MW-9

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 06/10/09 13:45
Date Received: 06/11/09

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:46 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.029 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 11:00 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:46 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:29 | EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Lab ID: L0907670-05 Date Collected: 06/10/09 14:05

Client ID: Date Received: 06/11/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:52 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.066 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 11:02 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:52 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:32 | EPA 3005A | 60,6010B | Al |



06/10/09 14:12

See Narrative

06/11/09

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L0907670-06

Client ID: RIZ-3

Sample Location: WALPOLE, MA Field Prep:

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:57 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.013 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 11:04 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 01:57 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:42 | EPA 3005A | 60,6010B | Al |
| | | | | | | | | | | |



06/10/09 14:26

Project Name: Lab Number: WALPOLE PARK SOUTH L0907670

Project Number: 12700053 Report Date: 06/17/09

SAMPLE RESULTS

Date Collected:

Lab ID: L0907670-07

Client ID: MW-2

Date Received: 06/11/09 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|------------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | tals - Wes | tborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 02:03 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.070 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 11:05 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 02:03 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:45 | EPA 3005A | 60,6010B | AI |



06/10/09 14:41

06/11/09

Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L0907670-08

Client ID: RIZ-9

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 02:09 | EPA 3005A | 64,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.015 | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 11:11 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 06/12/09 11:00 | 06/17/09 02:09 | EPA 3005A | 64,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:48 | EPA 3005A | 60,6010B | Al |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 Report Date: 06/17/09

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------|----------------------|----------|-----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals | - Westborough Lab fo | r sample | (s): 01-C | 8 Batch: | WG366599-1 | | | |
| Arsenic, Dissolved | ND | mg/l | 0.005 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Barium, Dissolved | ND | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Beryllium, Dissolved | ND | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Cadmium, Dissolved | ND | mg/l | 0.004 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Chromium, Dissolved | ND | mg/l | 0.01 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Lead, Dissolved | ND | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Nickel, Dissolved | ND | mg/l | 0.025 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Selenium, Dissolved | ND | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Silver, Dissolved | ND | mg/l | 0.007 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Vanadium, Dissolved | ND | mg/l | 0.010 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |
| Zinc, Dissolved | ND | mg/l | 0.050 | 1 | 06/12/09 11:00 | 06/15/09 16:07 | 60,6010B | AI |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|------------------------|--------------------|----------|----------|--------------------|------------------|------------------|----------------------|----|
| MCP Dissolved Metals - | Westborough Lab fo | r sample | (s): 01- | 08 Batch: | WG366913-1 | | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 06/15/09 17:15 | 06/16/09 10:49 | 64,7470A | EZ |

Prep Information

Digestion Method: EPA 7470A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------|----------------------|----------|----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals | - Westborough Lab fo | r sample | (s): 01- | 08 Batch: | WG367081-1 | | | |
| Antimony, Dissolved | ND | mg/l | 0.0005 | 1 | 06/12/09 11:00 | 06/17/09 00:42 | 64,6020A | ВМ |
| Thallium, Dissolved | ND | mg/l | 0.0005 | 1 | 06/12/09 11:00 | 06/17/09 00:42 | 64,6020A | ВМ |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 **Report Date:** 06/17/09

Method Blank Analysis Batch Quality Control

Prep Information

Digestion Method: EPA 3005A



Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053

Lab Number: L0907670

| arameter | LCS %Recovery | % | LCSD Recovery | %Recovery Limits | RPD | RPD Limits |
|--|-----------------------|-------|------------------|-----------------------|-----|------------|
| MCP Dissolved Metals - Westborough Lab | Associated sample(s): | 01-08 | Batch: \ | WG366599-2 WG366599-3 | | |
| Arsenic, Dissolved | 113 | | 114 | 80-120 | 1 | 20 |
| Barium, Dissolved | 106 | | 106 | 80-120 | 0 | 20 |
| Beryllium, Dissolved | 107 | | 106 | 80-120 | 1 | 20 |
| Cadmium, Dissolved | 116 | | 115 | 80-120 | 1 | 20 |
| Chromium, Dissolved | 105 | | 105 | 80-120 | 0 | 20 |
| Lead, Dissolved | 110 | | 111 | 80-120 | 1 | 20 |
| Nickel, Dissolved | 105 | | 105 | 80-120 | 0 | 20 |
| Selenium, Dissolved | 113 | | 117 | 80-120 | 3 | 20 |
| Silver, Dissolved | 110 | | 109 | 80-120 | 1 | 20 |
| Vanadium, Dissolved | 106 | | 105 | 80-120 | 1 | 20 |
| Zinc, Dissolved | 110 | | 111 | 80-120 | 1 | 20 |
| CP Dissolved Metals - Westborough Lab | Associated sample(s): | 01-08 | Batch: \ | WG366913-2 WG366913-3 | | |
| Mercury, Dissolved | 98 | | 93 | 80-120 | 5 | 20 |
| ICP Dissolved Metals - Westborough Lab | Associated sample(s): | 01-08 | Batch: \ | WG367081-2 WG367081-3 | | |
| Antimony, Dissolved | 106 | | 106 | 80-120 | 2 | 20 |
| Thallium, Dissolved | 96 | | 96 | 80-120 | 1 | 20 |



Lab Number: L0907670

Project Name: WALPOLE PARK SOUTH

Project Number: 12700053 Report Date: 06/17/09

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal A Absent

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|----------------------------------|--------|-----|------|------|--------|--|
| L0907670-01A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-01B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-01C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-TA70S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-02A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-02B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-02C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-03A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-03B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-03C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-04A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-04B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |



Project Name: WALPOLE PARK SOUTH

Lab Number: L0907670 Project Number: 12700053 **Report Date:** 06/17/09

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|----------------------------------|--------|-----|------|------|--------|--|
| L0907670-04C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-TL-6020S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-05A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-05B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-05C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-TL-6020S(180),MCP-TL-6020S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-06A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-06B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-06C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-TL-6020S(180),MCP-PB-6010S(180),MCP-PB-6010S(180),MCP-SN-6010S(180),MCP-SN-6010S(180),MCP-V-6010S(180) |
| L0907670-07A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-07B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-07C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-T470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0907670-08A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |
| L0907670-08B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 3 | Υ | Absent | 524.2(14) |



Project Name: WALPOLE PARK SOUTH Lab Number: L0907670

Project Number: 12700053 Report Date: 06/17/09

Container Information

| Container ID | Container Type | Cooler | рН | Temp | Pres | Seal | Analysis |
|--------------|------------------------------|--------|----|------|------|--------|--|
| L0907670-08C | Plastic 500ml HNO3 preserved | A | <2 | 3 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-T470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-R5-6010S(180),MCP-V-6010S(180) |



Project Name:WALPOLE PARK SOUTHLab Number:L0907670Project Number:12700053Report Date:06/17/09

GLOSSARY

Acronyms

EPA · Environmental Protection Agency.

LCS Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD · Laboratory Control Sample Duplicate: Refer to LCS.

MS • Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD · Matrix Spike Sample Duplicate: Refer to MS.

NA · Not Applicable.

NC · Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

ND · Not detected at the reported detection limit for the sample.

NI · Not Ignitable.

RDL • Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- * The batch duplicate RPD exceeds the acceptance criteria. This flag is not applicable when the sample concentrations are less than 5x the RDL. (Metals only.)
- A Spectra identified as "Aldol Condensation Product".
- **B** The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E -Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- N The matrix spike recovery exceeds the acceptance criteria. This flag is not applicable when the sample concentration is greater than 4x the spike added. (Metals only.)
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).

Report Format: Data Usability Report



Project Name:WALPOLE PARK SOUTHLab Number:L0907670Project Number:12700053Report Date:06/17/09

REFERENCES

Methods for the Determination of Organic Compounds in Drinking Water - Supplement II. EPA/600/R-92/129, August 1992.

- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). May 2004.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised June 17, 2009 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. NELAP Accredited Solid Waste/Soil.

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate. Organic Parameters: Haloacetic Acids, Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB).) Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Calcium Hardness, Silica, Sulfate, Sulfide, Ammonia, Kieldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics.) Solid Waste/Soil (Inorganic Parameters: Lead in Paint, pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), Reactivity. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3.3'-Dichlorobenzidine, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9221E, 9222B, 9222D, 9223B, EPA 150.1, 180.1, 300.0, 353.2, SM2130B, 2320B, 4500Cl-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B,4500NO3-F, EPA 200.7, EPA 200.8, 245.1. Organic Parameters: 504.1, 524.2, SM 6251B.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, Lachat 10-107-06-1-B, SM2320B, 2340B, 2510B, 2540C, 2540D, 426C, 4500Cl-D, 4500Cl-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-G, 4500NH3-H, 4500NO3-F, 4500P-B.5, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water

Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl)

(EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Nitrite-N, Fluoride, Sulfate)

353.2 for: Nitrate-N, Nitrite-N; SM4500NO3-F, 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, 2320B, SM2540C, EPA 150.1, SM4500H-B.

Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics)

(504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), SM6251B, 314.0.

Non-Potable Water

Inorganic Parameters:, (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn) (EPA 200.7 for: Al,Sb,As,Be,Cd,Cr,Co,Cu,Fe,Pb,Mn,Mo,Ni,Se,Ag,Sr,Tl,Ti,V,Zn,Ca,Mg,Na,K) 245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2540B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Nitrate-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-B,C-Titr, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B,E, 5220D, EPA 410.4, SM 5210B, 5310C, 4500CN-CE, 2540D, 4500CL-D, EPA 1664, SM14 510AC, EPA 420.1

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics)

(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCB-Water)

600/4-81-045-PCB-Oil

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water

OQA-QAM-025 Rev.7.)

Microbiology Parameters: SM9215B; MF-SM9222B; ENZ. SUB. SM9223; EC-SM9221E; MF-SM9222D; ENZ. SUB. SM9223;

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM6215B, 9222B, 9223B Colilert, EPA 200.7, 200.8, 245.2, 110.2, 120.1, 150.1, 300.0, 325.2, 314.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 331.0. Organic Parameters: 504.1, 524.2, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 150.1, 300.0, 305.1, 310.1, 325.2, 340.2, 350.1, 350.2, 351.1, 353.2, 354.1, 365.2, 375.4, 376.2, 405.1, 415.1, 420.1, 425.1, 1664A, SW-846 9010, 9030, 9040B, EPA 160.1, 160.2, 160.3, SM426C, SM2310B, 2540B, 2540D, 4500H+B, 4500NH3-H, 4500NH3-E, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 2320B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-117-07-1-B, LACHAT 10-107-06-1-B, LACHAT 10-107-04-1-C, LACHAT 10-107-04-1-J, LACHAT 10-117-07-1-A, SM4500CL-E, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3005A, 3015A, 3510C, 5030B, 8021B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 7.3.3.2, 7.3.4.2, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040, 9045C, 9050C, 1311, 3005A, 3050B, 3051A. Organic Parameters: SW-846 3540C, 3545, 3580A, 5030B, 5035, 8021B, 8260B, 8270C, 8330, 8151A, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 2540C, 2320B, 314.0, 331.0, 110.2, SM2120B, 2510B, 5310C, EPA 150.1, SM4500H-B, EPA 200.8, 245.2. Organic Parameters: 504.1, SM6251B, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.1, SM5220D, 4500Cl-D, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, SM9221CE, 9222D, 9221B, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.2/.1, SM5210B, SW-846 3015, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 3510C, EPA 608, 624, 625, SW-846 5030B, 8021B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7.) Solid & Chemical Materials (Inorganic Parameters: SW-846 9040B, 3005A, 6010B, 7196A, 5030B, 9010B, 9030B, 1030, 1311, 3050B, 3051, 7471A, 9014, 9012A, 9045C, 9050A, 9065. Organic Parameters: SW-846 8021B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 1311, 1312, 3540C, 3545, 3550B, 3580A, 5035L, 5035H, NJ

New York Department of Health Certificate/Lab ID: 11148. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 8215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 314.0, 331.0, SM2320B, EPA 300.0, 325.2, 110.2, SM2120B, 4500CN-E, 4500F-C, EPA 150.1, SM4500H-B, 4500NO3-F, 2540C, EPA 120.1, SM 2510B. Organic Parameters: EPA 524.2, 504.1, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, EPA 405.1, SM5210B, EPA 410.4, SM5220D, EPA 305.1, SM2310B-4a, EPA 310.1, SM2320B, EPA 200.7, 300.0, 325.2, LACHAT 10-117-07-1A or B, SM4500Cl-E, EPA 340.2, SM4500F-C, EPA 375.4, SM15 426C, EPA 350.1, 350.2, LACHAT 10-107-06-1-B, SM4500NH3-H, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-041-C, SM4500-NO30F, EPA 354.1, SM4500-NO2-B, EPA 365.2, SM4500P-E, EPA 160.3, EPA 160.1, SM2540C, EPA 160.2, SM2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, S\M3500Cr-D, EPA 245.1, 245.2, 7470A, 110.2, SM2120B, 335.2, LACHAT 10-204-00-1-A, EPA 150.1, 9040B, SM4500-HB, EPA 1664A, EPA 415.1, SM5310C, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, EPA 376.2, SM4500S-D, EPA 425.1, SM5540C, EPA 3005A, 3015. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, 8021B, EPA 3510C, 5030B, 9010B, 9030B.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 9040B, 9045C, 1010, 1030, SW-846 Ch 7 Sec 7.3, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 3005A, 3050B, 3051, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8151A, 8330, 8082, 8021B, 3540C, 3545, 3580, 5030B, 5035.)

Analytical Services Protocol: CLP Volatile Organics, CLP Inorganics, CLP PCB/Pesticides.

Rhode Island Department of Health Certificate/Lab ID: LAO00065. NELAP Accredited via NY-DOH.

Refer to MA-DEP Certificate for Potable and Non-Potable Water.

Refer to NY-DOH Certificate for Potable and Non-Potable Water.

Pennsylvania Department of Environmental Protection Certificate/Lab ID: 68-03671. *NELAP Accredited. Non-Potable Water* (Organic Parameters: EPA 3510C, 625, 608, 8081A, 8082, 8151A, 8270C, 8330) *Solid & Hazardous Waste* (Inorganic Parameters: EPA 1010, 1030, 1311, 3050B, 3051, 6010B, EPA 7.3.3.2, EPA 7.3.4.2, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065. Organic Parameters: 3540C, 3545, 3580A, 5035, 8021B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

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| /, | Due Date: 6/8/69 Time: | ☐ These samples have been Previously analyzed by Alpha |
| | i | Email Jan. Cannan & telately. Con |
| PPROVED) | Rush (ONLY IF PRE-APPROVED) | Fax: 50 93 7001 |
| ANALYSIS | Turn-Around Time | Phone: 508 905 2039 |
| | ALPHA Quote #: | Famingham Mit |
| | Project Manager: Nay Jurson | is: |
| MA MCF | Project #: 1270005 3 | Client Tetlatak Raz- |
| m Criteria | Project Location: Walpak, NVT | Client Information |
| Regulatory Requirements/Report Limits | WINDLE MEK JOHN | |
| ADEX Addi Deliverables | Project Name: | Westborough, MA Mansfield, MA TEL: 508-898-9220 TEL: 508-822-9300 |
| M EMAIL | | 0 |
| Report Information Data Deliverables Billing Information | | NEPHA |
| Date Rec'd in Lab Co [1/1/69 ALPHA Job #: | CHAIN OF CUSTODY PAGE 1 OF 1 | CHAIN OF |



ANALYTICAL REPORT

Lab Number: L1003740

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: Ray Johnson

Project Name: WALPOLE PK SOUTH

Project Number: 12700058 Report Date: 03/19/10

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name:WALPOLE PK SOUTHLab Number:L1003740

Project Number: 12700058 **Report Date:** 03/19/10

| Alpha Sample ID | Client ID | Sample Location | Collection Date/Time |
|--------------------|-----------|--------------------|-------------------------|
| L1003740-01 | MW-9 | WALPOLE, MA | 12/21/09 10:12 |
| L1003740-02 | RIZ-8 | WALPOLE, MA | 12/21/09 11:05 |
| L1003740-03 | GHC-6 | WALPOLE, MA | 12/28/09 08:25 |
| L1003740-04 | RIZ-3 | WALPOLE, MA | 12/28/09 09:03 |
| L1003740-05 | RIZ-9 | WALPOLE, MA | 12/28/09 10:00 |
| L1003740-06 | RIZ-10 | WALPOLE, MA | 12/28/09 10:43 |
| L1003740-07 | MW-3 | WALPOLE, MA | 12/28/09 11:50 |

Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: 12700058 Report Date: 03/19/10

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | NO |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L1003740

Lab Number:

Project Name: WALPOLE PK SOUTH

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

For additional information, please contact Client Services at 800-624-9220.

MCP Related Narratives

Sample Receipt

The samples were Field Filtered.

Metals

L1003740-01 through -07 have elevated detection limits due to the dilutions required by the high concentrations of non-target analytes. The requested reporting limits were achieved.

In reference to question F:

All samples were analyzed for a subset of MCP elements per the Chain of Custody.

Unabeth & Simuro

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative

ANALYTICA

Date: 03/19/10

METALS



Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

Lab ID: L1003740-01 Date Collected: 12/21/09 10:12

Client ID: MW-9 Date Received: 12/29/09

WALPOLE, MA Field Prep: Sample Location: See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst

MCP Dissolved Metals - Westborough Lab ND 64,6020A TD Lead, Dissolved mg/l 0.0020 4 12/30/09 09:40 12/30/09 18:56 EPA 3005A



Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

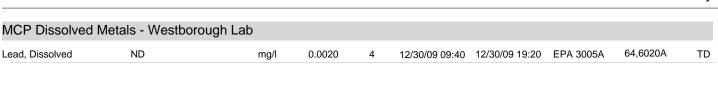
Lab ID: L1003740-02 Date Collected: 12/21/09 11:05

Client ID: RIZ-8 Date Received: 12/29/09

WALPOLE, MA Field Prep: Sample Location: See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst





Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: 12700058 **Report Date:** 03/19/10

SAMPLE RESULTS

Lab ID: L1003740-03 Date Collected: 12/28/09 08:25

Client ID: Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

Dilution Date Drep Analytical
Parameter Result Qualifier Units RDL Factor Prepared Analyzed Method Method Analyst

MCP Dissolved Metals - Westborough Lab

Lead, Dissolved ND mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:26 EPA 3005A 64,6020A TD



Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

Lab ID: L1003740-04 Date Collected: 12/28/09 09:03

Client ID: RIZ-3 Date Received: 12/29/09

Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst

MCP Dissolved Metals - Westborough Lab ND 64,6020A TD Lead, Dissolved mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:32 EPA 3005A



Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

Lab ID: L1003740-05 Date Collected: 12/28/09 10:00

RIZ-9 Client ID: Date Received: 12/29/09

WALPOLE, MA Field Prep: Sample Location: See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst

MCP Dissolved Metals - Westborough Lab ND 64,6020A TD Lead, Dissolved mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:38 EPA 3005A



Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

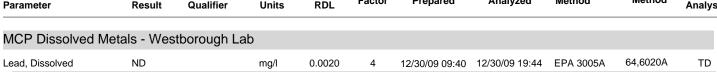
Lab ID: L1003740-06 Date Collected: 12/28/09 10:43

Client ID: RIZ-10 Date Received: 12/29/09

Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst





Project Name: WALPOLE PK SOUTH Lab Number: L1003740

Project Number: Report Date: 12700058 03/19/10

SAMPLE RESULTS

Lab ID: L1003740-07 Date Collected: 12/28/09 11:50

Client ID: MW-3 Date Received: 12/29/09

Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water

Analytical Method Dilution Date Date Prep Factor Prepared Analyzed Method Parameter Result Qualifier Units RDL Analyst

MCP Dissolved Metals - Westborough Lab ND 64,6020A TD Lead, Dissolved mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:50 EPA 3005A



Project Name: WALPOLE PK SOUTH **Lab Number:** L1003740

Project Number: 12700058 **Report Date:** 03/19/10

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|---------------------|-------------------------|----------|-----------|--------------------|------------------|------------------|----------------------|----|
| MCP Dissolved Metal | ls - Westborough Lab fo | r sample | e(s): 01- | 07 Batch: | WG404525-1 | | | |
| Lead, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/09 09:40 | 12/30/09 18:32 | 64,6020A | TD |

Prep Information

Digestion Method: EPA 3005A



Lab Control Sample Analysis

Batch Quality Control

Lab Number: L1003740

03/19/10

Project Number: 12700058 Report Date:

LCS **LCSD** %Recovery %Recovery %Recovery Limits **RPD Limits Parameter** Qual RPD Qual Qual MCP Dissolved Metals - Westborough Lab Associated sample(s): 01-07 Batch: WG404525-2 WG404525-3 Lead, Dissolved 103 101 80-120 2 20



Project Name:

WALPOLE PK SOUTH

Project Name:WALPOLE PK SOUTHLab Number:L1003740Project Number:12700058Report Date:03/19/10

GLOSSARY

Acronyms

EPA · Environmental Protection Agency.

LCS Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD · Laboratory Control Sample Duplicate: Refer to LCS.

MS • Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD · Matrix Spike Sample Duplicate: Refer to MS.

NA · Not Applicable.

NC · Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

NI · Not Ignitable.

RDL - Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A Spectra identified as "Aldol Condensation Product".
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- H -The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- ${f P}$ The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RDL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- J Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- ND Not detected at the reported detection limit (RDL) for the sample.

Report Format: Data Usability Report



Project Name:WALPOLE PK SOUTHLab Number:L1003740Project Number:12700058Report Date:03/19/10

REFERENCES

Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised March 16, 2010 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. NELAP Accredited Solid Waste/Soil.

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate. Organic Parameters: Haloacetic Acids, Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB).)

Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Calcium Hardness, Silica, Sulfate, Sulfide, Ammonia, Kjeldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH.) Solid Waste/Soil (Inorganic Parameters: Lead in Paint, pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), Reactivity. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH, Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3.3'-Dichlorobenzidine, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9221E, 9222B, 9222D, 9223B, EPA 180.1, 300.0, 353.2, SM2130B, 2320B, 4500Cl-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B,4500NO3-F, EPA 200.7, EPA 200.8, 245.1. Organic Parameters: 504.1, 524.2, SM 6251B.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, Lachat 10-107-06-1-B, SM2320B, 2340B, 2510B, 2540C, 2540D, 426C, 4500Cl-D, 4500Cl-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500P-B.5, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water

Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl)

(EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Fluoride, Sulfate)

353.2 for: Nitrate-N, Nitrite-N; SM4500NO3-F, 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, 2320B, SM2540C. SM4500H-B.

Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics)

(504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), 314.0, 332.

Microbiology Parameters: SM9215B; ENZ. SUB. SM9223; MF-SM9222D

Non-Potable Water

Inorganic Parameters:, (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn)

(EPA 200.7 for: Al,Sb,As,Be,Cd,Cr,Co,Cu,Fe,Pb,Mn,Mo,Ni,Se,Aq,Sr,Ti,Tl, V,Zn,Ca,Mq,Na,K)

245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2540B, 2340B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Ammonia-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-B,C-Titr, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B,E, 5220D, EPA 410.4, SM 5210B, 5310C, 4500CL-D, EPA 1664, SM14 510AC, EPA 420, SM4500-CN-CE, SM2540D.

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics)

(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCBs-Water), EPA 625 for SVOC Acid Extractables and SVOC Base/Neutral Extractables, 600/4-81-045-PCB-Oil

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM6215B, 9222B, 9223B Colilert, EPA 200.7, 200.8, 245.2, 120.1, 300.0, 314.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 331.0. Organic Parameters: 504.1, 524.2, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 300.0, 350.1, 351.1, 353.2, 420.1, 1664A, SW-846 9010, 9030, 9040B, SM426C, SM2310B, 2540B, 2540D, 4500H+B, 4500NH3-H, 4500NH3-E, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 2320B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-117-07-1-B, LACHAT 10-107-06-1-B, LACHAT 10-107-04-1-J, LACHAT 10-117-07-1-A, SM4500CL-E, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3005A, 3015A, 3510C, 5030B, 8021B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 7.3.3.2, 7.3.4.2, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040, 9045C, 9050C, 1311, 3005A, 3050B, 3051A. Organic Parameters: SW-846 3540C, 3545, 3580A, 5030B, 5035, 8021B, 8260B, 8270C, 8330, 8151A, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 2540C, 2320B, 314.0, SM2120B, 2510B, 5310C, SM4500H-B, EPA 200.8, 245.2. Organic Parameters: 504.1, SM6251B, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.4, SM5220D, 4500Cl-D, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, SM9221CE, 9222D, 9221B, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.1, SM5210B, SW-846 3015, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 3510C, EPA 608, 624, 625, SW-846 5030B, 8021B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 9040B, 3005A, 6010B, 7196A, 5030B, 9010B, 9030B, 1030, 1311, 3050B, 3051, 7471A, 9014, 9012A, 9045C, 9050A, 9065. Organic Parameters: SW-846 8021B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 1311, 1312, 3540C, 3545, 3550B, 3580A, 5035L, 5035H, NJ OQA-QAM-025 Rev.7.)

New York Department of Health Certificate/Lab ID: 11148. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 9215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 314.0, 332.0, SM2320B, EPA 300.0, SM2120B, 4500CN-E, 4500F-C, 4500H-B, 4500NO3-F, 2540C, EPA 120.1, SM 2510B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, 5210B, EPA 410.4, SM5220D, 2310B-4a, 2320B, EPA 200.7, 300.0, LACHAT 10-117-07-1A or B, SM4500Cl-E, 4500F-C, SM15 426C, EPA 350.1, LACHAT 10-107-06-1-B, SM4500NH3-H, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-041-C, SM4500-NO3-F, 4500-NO2-B, 4500P-E, 2540C, 2540B, 2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, S\M3500Cr-D, EPA 245.1, 245.2, 7470A, SM2120B, SM4500-CN-E LACHAT 10-204-00-1-A, EPA 9040B, SM4500-HB, EPA 1664A, SM5310C, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, SM4500S-D, SM5540C, EPA 3005A, 3015. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, EPA 3510C, 5030B, 9010B, 9030B.)

Solid & Hazardous Waste (Inorganic Parameters: 1010, 1030, SW-846 Ch 7 Sec 7.3, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 1312, 3005A, 3050B, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8151A, 8330, 8082, 3540C, 3545, 3546, 3580, 5030B, 5035.)

North Carolina Department of the Environment and Natural Resources <u>Certificate/Lab ID</u>: 666. <u>Organic Parameters</u>: MA-EPH, MA-VPH.

Pennsylvania Department of Environmental Protection Certificate/Lab ID: 68-03671. *NELAP Accredited. Non-Potable Water* (Organic Parameters: EPA 3510C, 5030B, 625, 624. 608, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1010, 1030, 1311, 3050B, 3051, 6010B, EPA 7.3.3.2, EPA 7.3.4.2, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065. Organic Parameters: 3540C, 3545, 3580A, 5035, 8021B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Rhode Island Department of Health <u>Certificate/Lab ID</u>: LAO00065. *NELAP Accredited via NY-DOH*. Refer to MA-DEP Certificate for Potable and Non-Potable Water. Refer to NY-DOH Certificate for Potable and Non-Potable Water.

Texas Commisson on Environmental Quality Certificate/Lab ID: T104704476-09-1. **NELAP Accredited.** Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664, 200.7, 200.8, 245.1, 245.2, 300.0, 350.1, 351.1, 353.2, 376.2, 410.4, 420.1, 6010, 6020, 7196, 7470, 9040, SM 2120B, 2310B, 2320B, 2510B, 2540B, 2540B, 2540D, 426C, 4500CL-E, 4500CN-E, 4500F-C, 4500H+B, 4500NH3-H, 4500NO2B, 4500P-E, 4500 S2⁻ D, 510C, 5210B, 5220D, 5310C, 5540C. Organic Parameters: EPA 608, 624, 625, 8081, 8082, 8151, 8260, 8270, 8330.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 9012, 9014, 9040, 9045, 9050, 9065.)

Utah Department of Health <u>Certificate/Lab ID</u>: AAMA. **NELAP Accredited. Non-Potable Water** (Inorganic Parameters: Chloride EPA 300.0)

Department of Defense Certificate/Lab ID: L2217.

Drinking Water (Inorganic Parameters: SM 4500H-B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: EPA 200.7, 200.8, 6010B, 6020, 245.1, 245.2, 7470A, 9040B, 300.0, 9251, 9038, 350.1, 353.2, 351.1, 314, 120.1, 9050A, 410.4, 9060, 1664, 420.1, LACHAT 10-107-06-1-B, SM 4500CN-E, 4500H-B, 4500CL-E, 4500F-BC, 4500SO4-E, 426C, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500NO2-B, 4500Norg-C, 4500PE, 2510B, 5540C, 5220D, 5310C, 2540B, 2540C, 2540D, 510C, 4500S2-AD, 3005A, 3015, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8330, 625, 8082, 8151A, 8081A, 3510C, 5030B.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 200.7, 6010B, 7471A, 9040B, 9045C, 9065, 420.1, 9012A, 6860, 1311, 1312, 3050B, 9030B, 3051, 9010B, 3540C, SM 510ABC, 4500CN-CE, 2540G, SW-846 7.3, Organic Parameters: EPA 8260B, 8270C, 8330, 8082, 8081A, 8151A, 3545, 3546, 3580, 5035.)

Analytes Not Accredited by NELAP

Certification is not available by NELAP for the following analytes: **EPA 8260B:** Freon-113, 1,2,4,5-Tetramethylbenzene, 4-Ethyltoluene. **EPA 8330A:** PETN, Picric Acid, Nitroglycerine, 2,6-DANT, 2,4-DANT. **EPA 8270C:** Methyl naphthalene, Dimethyl naphthalene, Total Methylnapthalenes, Total Dimethylnaphthalenes, 1,4-Diphenylhydrazine (Azobenzene). **EPA 625:** 4-Chloroaniline. **EPA 350.1** for Ammonia in a Soil matrix.

| PLEASE ANSWER QUESTIONS ABOVE! Container Type IS YOUR PROJECT MA MCP or CT RCP? Common of the 14-001-07) 8 6HC-6 10/28/4 0725 10500 S R1Z-2 10600 1043 1 1043 1 10600 1043 1 10600 1043 1 10600 106 | ments/Comments/Detection L ments/Comments/Detection L Coll ample ID Coll Date | CHAIN OF CUSTODY PAGE 1 OF 1 AMANSFIELD, MA TEL: 508-822-2000 FAX: 508-822-3288 Client Information Client Felix Felt 17220 Project Location: Majale M. Saul Project Manager: Majale M. Saul Project M |
|--|--|--|--|
| Hcl // Please prints early legibly and com- pletely. Samples can not be legged in and furnation time clock will not start unit any ambiguities are resolved All samples submitted are seconed All samples submitted are subjective (72/74/04/125) Seere erse sub- | | A Samp | ALEPHA Job # CONFIDENCE PROTOCOL ACTUAL ACTUAL ACTUAL ACTUAL ACTUAL ACTUAL ACTUAL Criteria PRESUMPTIVE CERTAINTY - CT REASONABLE CONFIDENCE PROTOCOL DNo Are MCP Analytical Methods Required? SYNO Are CT RCP (Reasonable Confidence Protocols) Required? SYNO SAMPLE HANDLING |

.



ANALYTICAL REPORT

Lab Number: L0918777

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: lan Cannan

Project Name: WALPOLE PK SOUTH

Project Number: 12700058 Report Date: 01/05/10

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

| Alpha Sample ID | Client ID | Sample Location | Collection Date/Time |
|--------------------|------------|--------------------|-------------------------|
| L0918777-01 | MW-9 | WALPOLE, MA | 12/21/09 10:12 |
| L0918777-02 | RIZ-8 | WALPOLE, MA | 12/21/09 11:05 |
| L0918777-03 | GHC-6 | WALPOLE, MA | 12/28/09 08:25 |
| L0918777-04 | RIZ-3 | WALPOLE, MA | 12/28/09 09:03 |
| L0918777-05 | RIZ-9 | WALPOLE, MA | 12/28/09 10:00 |
| L0918777-06 | RIZ-10 | WALPOLE, MA | 12/28/09 10:43 |
| L0918777-07 | MW-3 | WALPOLE, MA | 12/28/09 11:50 |
| L0918777-08 | TRIP BLANK | WALPOLE, MA | 12/21/09 00:00 |

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 Report Date: 01/05/10

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | YES |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L0918777

Lab Number:

Project Name: WALPOLE PK SOUTH

Project Number: 12700058 **Report Date:** 01/05/10

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

| For addition | al information | nlease contact | Client Services | at 800-624-9220 |
|--------------|------------------------|------------------|-----------------|-----------------|
| oi auditioi | iai ii ii Ulii laiiUli | . Dicase cultact | CHELL DELVICES | al 000-024-3220 |

MCP Related Narratives

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

Metals

L0918777-01 through -07 have elevated detection limits for Antimony and Thallium due to the dilutions required by the high concentrations of non-target analytes. The requested reporting limits were achieved.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 01/05/10

Michelle M. Morris

ORGANICS



VOLATILES



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: 12/21/09 10:12

Client ID: MW-9 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 12/30/09 11:35

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|---------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborou | ıgh Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-01 12/21/09 10:12

Client ID: MW-9 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Facto |
|-------------------------------------|------------|-----------|-------|------|----------------|
| Volatile Organics by GC/MS - Westbo | orough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| o-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | 0.77 | | ug/l | 0.50 | 1 |

ND

ug/l



1

No Tentatively Identified Compounds

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: 12/21/09 10:12

Client ID: MW-9 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-02 12/21/09 11:05

Client ID: RIZ-8

Date Received: 12/29/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 12/30/09 12:12

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbor | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-02 Date Collected: 12/21/09 11:05

Client ID: RIZ-8 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Facto |
|---------------------------------------|----------|-----------|-------|------|----------------|
| Volatile Organics by GC/MS - Westbord | ough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-02 Date Collected: 12/21/09 11:05

Client ID: RIZ-8 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 | |
| 4-Bromofluorobenzene | 93 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-03 12/28/09 08:25

Client ID: GHC-6

Date Received: 12/29/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 12/30/09 12:49 Analytical Date:

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------|-----------|-------|------|------------------------|
| Volatile Organics by GC/MS - Westborough | n Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Date Collected: Lab ID: L0918777-03 12/28/09 08:25

Client ID: GHC-6 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Campio Location. While CLL, whi | | 1 101 | riola riop. | | |
|--|--------|-----------|-------------|------|------------------------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
| Volatile Organics by GC/MS - Westborough | Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| ,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| ,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| ,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| -Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| lexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| sopropylbenzene | ND | | ug/l | 0.50 | 1 |
| -lsopropyltoluene | ND | | ug/l | 0.50 | 1 |
| laphthalene | ND | | ug/l | 0.50 | 1 |
| -Propylbenzene | ND | | ug/l | 0.50 | 1 |
| ec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| ert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| ,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| ,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| ,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| ,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| romobenzene | ND | | ug/l | 0.50 | 1 |
| -Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| -Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| ,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| ,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| ,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| Footstively Identified Company | | | | | |
| Tentatively Identified Compounds | | | | | |
| | | | | | |

ND

ug/l



1

No Tentatively Identified Compounds

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-03 Date Collected: 12/28/09 08:25

Client ID: GHC-6 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-04 12/28/09 09:03

Client ID: RIZ-3

Date Received: 12/29/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 12/30/09 13:26

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | jh Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-04 Date Collected: 12/28/09 09:03

Client ID: RIZ-3 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| latile Organics by GC/MS - Westborough La Dichlorobenzene Dichlorobenzene ene ylene | Result b ND ND ND | Qualifier | Units ug/l | RDL | Dilution Factor |
|---|----------------------|-----------|---------------|------|-----------------|
| Dichlorobenzene Dichlorobenzene ene | ND ND | | ug/l | | |
| Dichlorobenzene ene | ND | | ug/l | | |
| Dichlorobenzene ene | ND | | ug/i | 0.50 | 1 |
| ene | | | | | 1 |
| | [311.1 | | ug/l | 0.50 | 1 |
| | | | ug/l | 0.50 | 1 |
| | ND | | ug/l | 0.50 | 1 |
| Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| mochloromethane | ND | | ug/l | 0.50 | 1 |
| utylbenzene | ND | | ug/l | 0.50 | 1 |
| nlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| achlorobutadiene | ND | | ug/l | 0.50 | 1 |
| ropylbenzene | ND | | ug/l | 0.50 | 1 |
| opropyltoluene | ND | | ug/l | 0.50 | 1 |
| hthalene | ND | | ug/l | 0.50 | 1 |
| ropylbenzene | ND | | ug/l | 0.50 | 1 |
| Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| mobenzene | ND | | ug/l | 0.50 | 1 |
| hlorotoluene | ND | | ug/l | 0.50 | 1 |
| hlorotoluene | ND | | ug/l | 0.50 | 1 |
| omomethane | ND | | ug/l | 0.50 | 1 |
| Dibromoethane | ND | | ug/l | 0.50 | 1 |
| Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| Dichloropropane | ND | | ug/l | 0.50 | 1 |
| hyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| | | | | | |
| | | | | | |
| ntatively Identified Compounds | | | | | |

ND

ug/l



1

No Tentatively Identified Compounds

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: 12/28/09 09:03

Client ID: RIZ-3 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-05 12/28/09 10:00

RIZ-9 Client ID:

Date Received: 12/29/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 12/30/09 14:02

Analyst: TT

| Volatile Organics by GC/MS - Westborough Lab Methylene chloride ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichklorofucorethane ND ug/l 0.50 1 1_2-Dichloromethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform <th>Parameter</th> <th>Result</th> <th>Qualifier</th> <th>Units</th> <th>RDL</th> <th>Dilution Factor</th> | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane | Volatile Organics by GC/MS - Westborou | gh Lab | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itaris-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bernache ND ug/l 0.50 1 Bernacher ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 </td <td>1,1-Dichloroethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromodichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromofichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Eenzene ND ug/l 0.50 1 Tolluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorotluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Chloro | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itrans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <td< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| ND | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ehylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND ug/l 0.50 1 | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-05 Date Collected: 12/28/09 10:00

Client ID: RIZ-9 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Campio Location. With CLL, With | | | 1 101 | a i iop. | Coo Harrain |
|---|--------|-----------|-------|----------|-----------------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
| Volatile Organics by GC/MS - Westboroug | jh Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| Tantativaly Identified Comments | | | | | |
| Tentatively Identified Compounds | | | | | |
| | | | | | |

ND

ug/l



1

No Tentatively Identified Compounds

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-05 Date Collected: 12/28/09 10:00

Client ID: RIZ-9 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 107 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: Report Date: 12700058 01/05/10

SAMPLE RESULTS

Lab ID: Date Collected: L0918777-06 12/28/09 10:43

Client ID: RIZ-10

Date Received: 12/29/09 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 12/30/09 14:39

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborough La | ıb | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-06 Date Collected: 12/28/09 10:43

Client ID: RIZ-10 Date Received: 12/29/09

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Ouripic Location. With OLL, Wi | ` | 1 101 | аттор. | Occ Harrati | |
|---------------------------------------|----------|-----------|--------|-------------|------------------------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
| Volatile Organics by GC/MS - Westbord | ough Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| sopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |
| Tentatively Identified Compounds | | | | | |
| No Tentatively Identified Compounds | ND | | ug/l | | 1 |
| · · | | | | | |

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-06 Date Collected: 12/28/09 10:43

Client ID: RIZ-10 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 104 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



12/29/09

See Narrative

Date Received:

Field Prep:

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-07 Date Collected: 12/28/09 11:50

Client ID: MW-3

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 12/30/09 15:16

Analyst: TT

| Volatile Organics by GC/MS - Westborough Lab Methylene chloride ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Inchipotopopane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Totrachloroethane ND ug/l 0.50 1 Totrachloroethane ND ug/l 0.50 1 Tichloroforomethane ND ug/l 0.50 1 1_2-Dichloromethane ND ug/l 0.50 1 1_2-Dichloropropene ND ug/l 0.50 1 1_2-Dichloropropene ND ug/l 0.50 1 1_2-Dichloropropene ND ug/l 0.50 1 1_2-Dichloro | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane | Volatile Organics by GC/MS - Westborou | gh Lab | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichlorofuloromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itaris-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bernache ND ug/l 0.50 1 Bernacher ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 </td <td>1,1-Dichloroethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromodichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromofichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Eenzene ND ug/l 0.50 1 Tolluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorotluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Chloro | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itrans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <td< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| ND | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ehylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND ug/l 0.50 1 | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-07 Date Collected: 12/28/09 11:50

Client ID: MW-3 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL **Dilution Factor** Parameter Qualifier Units Result Volatile Organics by GC/MS - Westborough Lab ND 1,3-Dichlorobenzene ug/l 0.50 1 ND 0.50 1 1,4-Dichlorobenzene ug/l ND 0.50 Styrene ug/l 1 o-Xylene ND ug/l 0.50 1 ND 0.50 1,1-Dichloropropene ug/l 1 ND 0.50 2,2-Dichloropropane ug/l 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l ND 0.50 1 1,2,3-Trichloropropane ug/l Bromochloromethane ND ug/l 0.50 1 ND 0.50 n-Butylbenzene ug/l 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND 0.50 ug/l 1 tert-Butylbenzene ND ug/l 0.50 1 ND 0.50 1 1,2,3-Trichlorobenzene ug/l ND 0.50 1 1,2,4-Trichlorobenzene ug/l 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND 0.50 1 ug/l Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND 0.50 1 ug/l ug/l ND p-Chlorotoluene 0.50 1 Dibromomethane ND ug/l 0.50 1

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |

ND

ND

ND

ND



1

1

1

1

0.50

0.50

0.50

0.50

ug/l

ug/l

ug/l

ug/l

1,2-Dibromoethane

1,3-Dichloropropane

Methyl tert butyl ether

1,2-Dibromo-3-chloropropane

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-07 Date Collected: 12/28/09 11:50

Client ID: MW-3 Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-08 Date Collected: 12/21/09 00:00

Client ID: TRIP BLANK Date Received: 12/29/09 Sample Location: WALPOLE, MA Field Prep: None

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 12/30/09 15:53

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--------------------------------------|----------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westbor | ough Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-08 Date Collected: 12/21/09 00:00

Client ID: TRIP BLANK Date Received: 12/29/09 Sample Location: WALPOLE, MA Field Prep: None

| Parameter | Result | Qualifier U | nits RDL | Dilution Facto |
|-------------------------------------|------------|-------------|----------|----------------|
| Volatile Organics by GC/MS - Westbo | orough Lab | | | |
| 1,3-Dichlorobenzene | ND | u | g/l 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | u | g/l 0.50 | 1 |
| Styrene | ND | u | g/l 0.50 | 1 |
| o-Xylene | ND | u | g/l 0.50 | 1 |
| 1,1-Dichloropropene | ND | u | g/l 0.50 | 1 |
| 2,2-Dichloropropane | ND | u | g/l 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | u | g/l 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | u | g/l 0.50 | 1 |
| Bromochloromethane | ND | u | g/l 0.50 | 1 |
| n-Butylbenzene | ND | u | g/l 0.50 | 1 |
| Dichlorodifluoromethane | ND | u | g/I 0.50 | 1 |
| Hexachlorobutadiene | ND | u | g/l 0.50 | 1 |
| sopropylbenzene | ND | u | g/I 0.50 | 1 |
| o-Isopropyltoluene | ND | u | g/l 0.50 | 1 |
| Naphthalene | ND | u | g/l 0.50 | 1 |
| n-Propylbenzene | ND | u | g/l 0.50 | 1 |
| sec-Butylbenzene | ND | u | g/I 0.50 | 1 |
| ert-Butylbenzene | ND | u | g/I 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | u | g/I 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | u | g/I 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | u | g/I 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | u | g/l 0.50 | 1 |
| Bromobenzene | ND | u | g/I 0.50 | 1 |
| o-Chlorotoluene | ND | u | g/I 0.50 | 1 |
| o-Chlorotoluene | ND | u | g/l 0.50 | 1 |
| Dibromomethane | ND | u | g/l 0.50 | 1 |
| 1,2-Dibromoethane | ND | u | g/l 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | u | g/I 0.50 | 1 |
| 1,3-Dichloropropane | ND | u | g/l 0.50 | 1 |
| Methyl tert butyl ether | ND | u | g/l 0.50 | 1 |

ND

ug/l



1

No Tentatively Identified Compounds

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-08 Date Collected: 12/21/09 00:00

Client ID: TRIP BLANK Date Received: 12/29/09
Sample Location: WALPOLE, MA Field Prep: None

Parameter Result Qualifier Units RDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 106 | | 80-120 | |
| 4-Bromofluorobenzene | 93 | | 80-120 | |



L0918777

Project Name: WALPOLE PK SOUTH Lab Number:

Project Number: 12700058 **Report Date:** 01/05/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 12/30/09 07:18

Analyst: TT

| arameter | Result | Qualifier | Units | RDL | |
|---------------------------|------------------|------------------|---------|--------|------------|
| olatile Organics by GC/MS | - Westborough La | b for sample(s): | : 01-08 | Batch: | WG395058-2 |
| Methylene chloride | ND | | ug/l | 0.8 | 50 |
| 1,1-Dichloroethane | ND | | ug/l | 0.9 | 50 |
| Chloroform | ND | | ug/l | 0.9 | 50 |
| Carbon tetrachloride | ND | | ug/l | 0.9 | 50 |
| 1,2-Dichloropropane | ND | | ug/l | 0.9 | 50 |
| Dibromochloromethane | ND | | ug/l | 0.9 | 50 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.9 | 50 |
| Tetrachloroethene | ND | | ug/l | 0.9 | 50 |
| Chlorobenzene | ND | | ug/l | 0.9 | 50 |
| Trichlorofluoromethane | ND | | ug/l | 0.9 | 50 |
| 1,2-Dichloroethane | ND | | ug/l | 0.9 | 50 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.9 | 50 |
| Bromodichloromethane | ND | | ug/l | 0.9 | 50 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.9 | 50 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.9 | 50 |
| Bromoform | ND | | ug/l | 0.9 | 50 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.9 | 50 |
| Benzene | ND | | ug/l | 0.9 | 50 |
| Toluene | ND | | ug/l | 0.9 | 50 |
| Ethylbenzene | ND | | ug/l | 0.9 | 50 |
| p/m-Xylene | ND | | ug/l | 0.9 | 50 |
| Chloromethane | ND | | ug/l | 0.9 | 50 |
| Bromomethane | ND | | ug/l | 0.9 | 50 |
| Vinyl chloride | ND | | ug/l | 0.9 | 50 |
| Chloroethane | ND | | ug/l | 0.9 | 50 |
| 1,1-Dichloroethene | ND | | ug/l | 0.9 | 50 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.9 | 50 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.9 | 50 |
| Trichloroethene | ND | | ug/l | 0.9 | 50 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.9 | 50 |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.9 | 50 |



L0918777

Project Name: WALPOLE PK SOUTH Lab Number:

Project Number: 12700058 **Report Date:** 01/05/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 12/30/09 07:18

Analyst: TT

| arameter | Result | Qualifier | Units | RDL | |
|-----------------------------|------------------|-----------------|-------|--------|------------|
| olatile Organics by GC/MS | - Westborough La | b for sample(s) | 01-08 | Batch: | WG395058-2 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.5 | 50 |
| Styrene | ND | | ug/l | 0.5 | 50 |
| o-Xylene | ND | | ug/l | 0.5 | 50 |
| 1,1-Dichloropropene | ND | | ug/l | 0.5 | 50 |
| 2,2-Dichloropropane | ND | | ug/l | 0.8 | 50 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.5 | 50 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.8 | 50 |
| Bromochloromethane | ND | | ug/l | 0.8 | 50 |
| n-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| Dichlorodifluoromethane | ND | | ug/l | 0.5 | 50 |
| Hexachlorobutadiene | ND | | ug/l | 0.5 | 50 |
| Isopropylbenzene | ND | | ug/l | 0.5 | 50 |
| p-Isopropyltoluene | ND | | ug/l | 0.5 | 50 |
| Naphthalene | ND | | ug/l | 0.5 | 50 |
| n-Propylbenzene | ND | | ug/l | 0.5 | 50 |
| sec-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| tert-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.8 | 50 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.5 | 50 |
| Bromobenzene | ND | | ug/l | 0.5 | 50 |
| o-Chlorotoluene | ND | | ug/l | 0.5 | 50 |
| p-Chlorotoluene | ND | | ug/l | 0.5 | 50 |
| Dibromomethane | ND | | ug/l | 0.5 | 50 |
| 1,2-Dibromoethane | ND | | ug/l | 0.8 | 50 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.5 | 50 |
| 1,3-Dichloropropane | ND | | ug/l | 0.8 | 50 |
| Methyl tert butyl ether | ND | | ug/l | 0.5 | 50 |
| | | | | | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 12/30/09 07:18

Analyst: TT

Parameter Result Qualifier Units RDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-08 Batch: WG395058-2

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | | 1 | Acceptance | |
|------------------------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | Criteria | |
| | | | | |
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 | |
| 4-Bromofluorobenzene | 95 | | 80-120 | |



Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| arameter | LCS %Recovery | Qual | | SD covery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|-------------------|-----------|-------|--------------|-----------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough I | _ab Associated sa | ample(s): | 01-08 | Batch: | WG395058- | 1 | | | |
| Methylene chloride | 96 | | | - | | 70-130 | - | | |
| 1,1-Dichloroethane | 103 | | | - | | 70-130 | - | | |
| Chloroform | 108 | | | - | | 70-130 | - | | |
| Carbon tetrachloride | 115 | | | - | | 70-130 | - | | |
| 1,2-Dichloropropane | 106 | | | - | | 70-130 | - | | |
| Dibromochloromethane | 111 | | | - | | 70-130 | - | | |
| 1,1,2-Trichloroethane | 107 | | | - | | 70-130 | - | | |
| Tetrachloroethene | 111 | | | - | | 70-130 | - | | |
| Chlorobenzene | 106 | | | - | | 70-130 | - | | |
| Trichlorofluoromethane | 115 | | | - | | 70-130 | - | | |
| 1,2-Dichloroethane | 118 | | | - | | 70-130 | - | | |
| 1,1,1-Trichloroethane | 116 | | | - | | 70-130 | - | | |
| Bromodichloromethane | 111 | | | - | | 70-130 | - | | |
| trans-1,3-Dichloropropene | 110 | | | - | | 70-130 | - | | |
| cis-1,3-Dichloropropene | 112 | | | - | | 70-130 | - | | |
| Bromoform | 119 | | | - | | 70-130 | - | | |
| 1,1,2,2-Tetrachloroethane | 105 | | | - | | 70-130 | - | | |
| Benzene | 104 | | | - | | 70-130 | - | | |
| Toluene | 104 | | | - | | 70-130 | - | | |
| Ethylbenzene | 105 | | | - | | 70-130 | - | | |
| p/m-Xylene | 106 | | | - | | 70-130 | - | | |

Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| ırameter | LCS %Recovery | Qual | | CSD covery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|------------|-------|---------------|-----------|---------------------|-----|------|------------|
| platile Organics by GC/MS - Westborough L | ab Associated s | sample(s): | 01-08 | Batch: | WG395058- | -1 | | | |
| Chloromethane | 104 | | | - | | 70-130 | - | | |
| Bromomethane | 91 | | | - | | 70-130 | - | | |
| Vinyl chloride | 101 | | | - | | 70-130 | - | | |
| Chloroethane | 102 | | | - | | 70-130 | - | | |
| 1,1-Dichloroethene | 100 | | | - | | 70-130 | - | | |
| trans-1,2-Dichloroethene | 101 | | | - | | 70-130 | - | | |
| cis-1,2-Dichloroethene | 103 | | | - | | 70-130 | - | | |
| Trichloroethene | 110 | | | - | | 70-130 | - | | |
| 1,2-Dichlorobenzene | 104 | | | - | | 70-130 | - | | |
| 1,3-Dichlorobenzene | 107 | | | - | | 70-130 | - | | |
| 1,4-Dichlorobenzene | 105 | | | - | | 70-130 | - | | |
| Styrene | 104 | | | - | | 70-130 | - | | |
| o-Xylene | 104 | | | - | | 70-130 | - | | |
| 1,1-Dichloropropene | 107 | | | - | | 70-130 | - | | |
| 2,2-Dichloropropane | 114 | | | - | | 70-130 | - | | |
| 1,1,1,2-Tetrachloroethane | 117 | | | - | | 70-130 | - | | |
| 1,2,3-Trichloropropane | 115 | | | - | | 70-130 | - | | |
| Bromochloromethane | 106 | | | - | | 70-130 | - | | |
| n-Butylbenzene | 99 | | | - | | 70-130 | - | | |
| Dichlorodifluoromethane | 115 | | | - | | 70-130 | - | | |
| Hexachlorobutadiene | 110 | | | - | | 70-130 | - | | |



Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| arameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---------------------------------------|-------------------|------------|-------------------|-----------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborou | gh Lab Associated | sample(s): | 01-08 Batch: | WG395058- | -1 | | | |
| Isopropylbenzene | 107 | | - | | 70-130 | - | | |
| p-Isopropyltoluene | 109 | | - | | 70-130 | - | | |
| Naphthalene | 72 | | - | | 70-130 | - | | |
| n-Propylbenzene | 103 | | - | | 70-130 | - | | |
| sec-Butylbenzene | 104 | | - | | 70-130 | - | | |
| tert-Butylbenzene | 109 | | - | | 70-130 | - | | |
| 1,2,3-Trichlorobenzene | 89 | | - | | 70-130 | - | | |
| 1,2,4-Trichlorobenzene | 91 | | - | | 70-130 | - | | |
| 1,2,4-Trimethylbenzene | 110 | | - | | 70-130 | - | | |
| 1,3,5-Trimethylbenzene | 110 | | - | | 70-130 | - | | |
| Bromobenzene | 109 | | - | | 70-130 | - | | |
| o-Chlorotoluene | 111 | | - | | 70-130 | - | | |
| p-Chlorotoluene | 107 | | - | | 70-130 | - | | |
| Dibromomethane | 101 | | - | | 70-130 | - | | |
| 1,2-Dibromoethane | 111 | | - | | 70-130 | - | | |
| 1,2-Dibromo-3-chloropropane | 99 | | - | | 70-130 | - | | |
| 1,3-Dichloropropane | 103 | | - | | 70-130 | - | | |
| Methyl tert butyl ether | 114 | | - | | 70-130 | - | | |

Lab Control Sample Analysis

Batch Quality Control

Lab Number: L0918777

Report Date: 01/05/10

LCS LCSD %Recovery

Parameter %Recovery Qual %Recovery Qual Limits RPD Qual RPD Limits

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 Batch: WG395058-1

WALPOLE PK SOUTH

12700058

| Surrogate | LCS %Recovery | Qual | LCSD %Recovery | Qual | Acceptance Criteria |
|------------------------|------------------|------|-------------------|------|------------------------|
| 1,2-Dichlorobenzene-d4 | 100 | | | | 80-120 |
| 4-Bromofluorobenzene | 101 | | | | 80-120 |



Project Name:

Project Number:

Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSI Qual Four | _ | Recovery y Qual Limits | RPE | RPD Qual Limits |
|---------------------------------------|------------------|-------------|--------------|-----------------|------------------|-------------|---------------------------|------|--------------------|
| olatile Organics by GC/MS - Sample | - Westborough | Lab Assoc | iated sample | (s): 01-08 Q | C Batch ID: WG | 395058-3 Q0 | C Sample: L091874 | 0-01 | Client ID: MS |
| Methylene chloride | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| Chloroform | ND | 4 | 4.5 | 112 | - | - | 70-130 | - | 20 |
| Carbon tetrachloride | ND | 4 | 4.9 | 123 | - | - | 70-130 | - | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| Dibromochloromethane | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 4.2 | 104 | - | - | 70-130 | - | 20 |
| Tetrachloroethene | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | 20 |
| Chlorobenzene | ND | 4 | 4.4 | 110 | - | - | 70-130 | - | 20 |
| Trichlorofluoromethane | ND | 4 | 5.0 | 124 | - | - | 70-130 | - | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.7 | 117 | - | - | 70-130 | - | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 5.2 | 130 | - | - | 70-130 | - | 20 |
| Bromodichloromethane | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| Bromoform | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | 20 |
| Benzene | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| Toluene | ND | 4 | 4.4 | 109 | - | - | 70-130 | - | 20 |
| Ethylbenzene | ND | 4 | 4.3 | 107 | - | - | 70-130 | - | 20 |
| p/m-Xylene | ND | 8 | 8.6 | 108 | - | - | 70-130 | - | 20 |
| | | | | | | | | | |

Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Qual Found | MSD %Recovery | Recovery Qual Limits | RPD | RPD Qual Limits |
|---------------------------------------|------------------|-------------|--------------|-----------------|-------------------|------------------|-------------------------|--------|--------------------|
| olatile Organics by GC/MS - cample | - Westborough | n Lab Assoc | iated sample | (s): 01-08 Q | C Batch ID: WG39 | 95058-3 QC | Sample: L091874 | 0-01 (| Client ID: MS |
| Chloromethane | ND | 4 | 3.6 | 91 | - | - | 70-130 | - | 20 |
| Bromomethane | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | 20 |
| Vinyl chloride | ND | 4 | 5.0 | 125 | - | - | 70-130 | - | 20 |
| Chloroethane | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.5 | 114 | - | - | 70-130 | - | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.2 | 106 | - | - | 70-130 | - | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 4.2 | 105 | - | - | 70-130 | - | 20 |
| Trichloroethene | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 4.0 | 100 | - | - | 70-130 | - | 20 |
| Styrene | ND | 4 | 3.9 | 97 | - | - | 70-130 | - | 20 |
| o-Xylene | ND | 4 | 4.1 | 104 | - | - | 70-130 | - | 20 |
| 1,1-Dichloropropene | ND | 4 | 4.4 | 111 | - | - | 70-130 | - | 20 |
| 2,2-Dichloropropane | ND | 4 | 4.8 | 120 | - | - | 70-130 | - | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 4.6 | 115 | - | - | 70-130 | - | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | 20 |
| Bromochloromethane | ND | 4 | 4.2 | 104 | - | - | 70-130 | - | 20 |
| n-Butylbenzene | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | 20 |
| Dichlorodifluoromethane | ND | 4 | 3.9 | 98 | - | - | 70-130 | - | 20 |
| Hexachlorobutadiene | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | 20 |
| | | | | | | | | | |

Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| rameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | Recovery Qual Limits | RPD | RPD Qual Limits |
|---------------------------------|------------------|-------------|-------------|-----------------|-----------|--------------|------------------|-------------------------|------|--------------------|
| olatile Organics by GC/MS ample | - Westborough | Lab Associ | ated sample | (s): 01-08 Q | C Batch I | D: WG395 | 5058-3 QC | Sample: L0918740 | 0-01 | Client ID: MS |
| Isopropylbenzene | ND | 4 | 3.8 | 96 | | - | - | 70-130 | - | 20 |
| p-Isopropyltoluene | ND | 4 | 4.3 | 107 | | - | - | 70-130 | - | 20 |
| Naphthalene | ND | 4 | 2.6 | 65 | Q | - | - | 70-130 | - | 20 |
| n-Propylbenzene | ND | 4 | 4.2 | 106 | | - | - | 70-130 | - | 20 |
| sec-Butylbenzene | ND | 4 | 4.3 | 108 | | - | - | 70-130 | - | 20 |
| tert-Butylbenzene | ND | 4 | 4.5 | 112 | | - | - | 70-130 | - | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.3 | 82 | | - | - | 70-130 | - | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.2 | 81 | | - | - | 70-130 | - | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 4.4 | 110 | | - | - | 70-130 | - | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 4.4 | 111 | | - | - | 70-130 | - | 20 |
| Bromobenzene | ND | 4 | 4.3 | 108 | | - | - | 70-130 | - | 20 |
| o-Chlorotoluene | ND | 4 | 4.4 | 110 | | - | - | 70-130 | - | 20 |
| p-Chlorotoluene | ND | 4 | 4.2 | 105 | | - | - | 70-130 | - | 20 |
| Dibromomethane | ND | 4 | 4.2 | 106 | | - | - | 70-130 | - | 20 |
| 1,2-Dibromoethane | ND | 4 | 4.4 | 109 | | - | - | 70-130 | - | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 3.7 | 93 | | - | - | 70-130 | - | 20 |
| 1,3-Dichloropropane | ND | 4 | 4.1 | 104 | | - | - | 70-130 | - | 20 |
| Methyl tert butyl ether | ND | 4 | 4.4 | 109 | <u> </u> | - | - | 70-130 | - | 20 |



Project Name: WALPOLE PK SOUTH

Project Number:

12700058

Lab Number:

L0918777

Report Date:

01/05/10

| | Native | MS | MS | MS | | MSD | MSD | | Recovery | | | RPD |
|-----------|--------|-------|-------|-----------|------|-------|-----------|------|----------|-----|------|--------|
| Parameter | Sample | Added | Found | %Recovery | Qual | Found | %Recovery | Qual | Limits | RPD | Qual | Limits |

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 QC Batch ID: WG395058-3 QC Sample: L0918740-01 Client ID: MS Sample

| | MS | MSD | Acceptance | |
|------------------------|----------------------|----------------------|------------|--|
| Surrogate | % Recovery Qualifier | % Recovery Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 98 | | 80-120 | |
| 4-Bromofluorobenzene | 102 | | 80-120 | |

Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number:

L0918777

Report Date:

| arameter | Native Sample | Duplicate Sample | Units | RPD | Qual RPD Limits |
|---|-----------------------------|----------------------|-------------|-------------|----------------------|
| platile Organics by GC/MS - Westborough Lab | Associated sample(s): 01-08 | QC Batch ID: WG39505 | 8-4 QC Samp | ole: L09187 | 83-02 Client ID: DUP |
| Methylene chloride | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethane | ND | ND | ug/l | NC | 20 |
| Chloroform | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene | ND | ND | ug/l | NC | 20 |
| Chlorobenzene | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| Bromoform | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| Benzene | ND | ND | ug/l | NC | 20 |
| Toluene | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number:

L0918777

Report Date:

| arameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|-----------------------------|------------------------|-------------|--------------|------------------|
| olatile Organics by GC/MS - Westborough Lab ample | Associated sample(s): 01-08 | QC Batch ID: WG395058- | 4 QC Sample | : L0918783-0 | 2 Client ID: DUP |
| Ethylbenzene | ND | ND | ug/l | NC | 20 |
| p/m-Xylene | ND | ND | ug/l | NC | 20 |
| Chloromethane | ND | ND | ug/l | NC | 20 |
| Bromomethane | ND | ND | ug/l | NC | 20 |
| Vinyl chloride | ND | ND | ug/l | NC | 20 |
| Chloroethane | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethene | ND | ND | ug/l | NC | 20 |
| trans-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| cis-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| Trichloroethene | ND | ND | ug/l | NC | 20 |
| 1,2-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,3-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,4-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| Styrene | ND | ND | ug/l | NC | 20 |
| o-Xylene | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloropropene | ND | ND | ug/l | NC | 20 |
| 2,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| 1,1,1,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichloropropane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number:

L0918777

Report Date:

| arameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------------|-------------------------|------------|-------------|----------------|
| olatile Organics by GC/MS - Westborough Lab ample | Associated sample(s): 01-08 | QC Batch ID: WG395058-4 | QC Sample: | L0918783-02 | Client ID: DUP |
| Bromochloromethane | ND | ND | ug/l | NC | 20 |
| n-Butylbenzene | ND | ND | ug/l | NC | 20 |
| Dichlorodifluoromethane | ND | ND | ug/l | NC | 20 |
| Hexachlorobutadiene | ND | ND | ug/l | NC | 20 |
| Isopropylbenzene | ND | ND | ug/l | NC | 20 |
| p-Isopropyltoluene | ND | ND | ug/l | NC | 20 |
| Naphthalene | ND | ND | ug/l | NC | 20 |
| n-Propylbenzene | ND | ND | ug/l | NC | 20 |
| sec-Butylbenzene | ND | ND | ug/l | NC | 20 |
| tert-Butylbenzene | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| 1,3,5-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| Bromobenzene | ND | ND | ug/l | NC | 20 |
| o-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| p-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| Dibromomethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dibromoethane | ND | ND | ug/l | NC | 20 |



WALPOLE PK SOUTH

Project Number: 12700058

Project Name:

Lab Number:

L0918777

Report Date:

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|---------------------------|-----------------------|-------------|-----------|----------------------|
| Volatile Organics by GC/MS - Westborough Lab Ass Sample | sociated sample(s): 01-08 | QC Batch ID: WG395058 | -4 QC Sampl | e: L09187 | 83-02 Client ID: DUP |
| 1,2-Dibromo-3-chloropropane | ND | ND | ug/l | NC | 20 |
| 1,3-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Methyl tert butyl ether | 1.4 | 1.4 | ug/l | 0 | 20 |

| | | | | | Acceptance | |
|------------------------|-----------|-----------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 102 | | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 93 | | 94 | | 80-120 | |



METALS



12/21/09 10:12

12/29/09

Project Name: WALPOLE PK SOUTH **Lab Number:** L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Dilution

Lab ID: L0918777-01

Client ID: MW-9

Sample Location: WALPOLE, MA

Matrix: Water

| Fie | ld Prep: | S | ee Narrative | 1 |
|----------|----------|--------|--------------|---------|
| Date | Date | Prep | Analytical | Analyst |
| Prepared | Analyzed | Method | Method | |

Date Collected:

Date Received:

| Parameter | Result | Qualifier | Units | RDL | Factor | Prepared | Analyzed | Method | Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------|----------------|----------------|-----------|----------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 18:56 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.056 | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:09 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 18:56 | EPA 3005A | 64,6020A | TD |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | 0.052 | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 17:57 | EPA 3005A | 60,6010B | AI |



12/21/09 11:05

12/29/09

Project Name: WALPOLE PK SOUTH **Lab Number:** L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L0918777-02

Client ID: RIZ-8

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:20 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | 0.005 | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.031 | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:11 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:20 | EPA 3005A | 64,6020A | TD |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 18:00 | EPA 3005A | 60,6010B | Al |



Project Name: Lab Number: WALPOLE PK SOUTH L0918777

Project Number: 12700058 Report Date: 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-03

Client ID: GHC-6

Sample Location: WALPOLE, MA

Matrix: Water Date Collected: 12/28/09 08:25

Date Received: 12/29/09

Field Prep: See Narrative

| Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|------------|--------------|---|--|---|--|--|---|--|---|
| tals - Wes | stborough La | ab | | | | | | | |
| ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:26 | EPA 3005A | 64,6020A | TD |
| ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| 0.039 | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:13 | EPA 7470A | 64,7470A | EZ |
| ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:26 | EPA 3005A | 64,6020A | TD |
| ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| ND | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 18:04 | EPA 3005A | 60,6010B | AI |
| | tals - Wes | tals - Westborough La ND ND 0.039 ND ND ND ND ND ND ND ND ND N | tals - Westborough Lab ND mg/l ND mg/l 0.039 mg/l ND mg/l | tals - Westborough Lab ND mg/l 0.0020 ND mg/l 0.005 0.039 mg/l 0.010 ND mg/l 0.004 ND mg/l 0.001 ND mg/l 0.010 ND mg/l 0.0002 ND mg/l 0.010 ND mg/l 0.007 ND mg/l 0.007 ND mg/l 0.0020 ND mg/l 0.0020 ND mg/l 0.0020 ND mg/l 0.0010 | Result Qualifier Units RDL Factor tals - Westborough Lab ND mg/l 0.0020 4 ND mg/l 0.005 1 0.039 mg/l 0.010 1 ND mg/l 0.004 1 ND mg/l 0.004 1 ND mg/l 0.010 1 ND mg/l 0.0002 1 ND mg/l 0.010 1 ND mg/l 0.007 1 ND mg/l 0.0020 4 ND mg/l 0.0020 4 ND mg/l 0.0020 4 ND mg/l 0.010 1 | Result Qualifier Units RDL Factor Prepared tals - Westborough Lab ND mg/l 0.0020 4 12/30/09 09:40 ND mg/l 0.005 1 01/04/10 11:15 0.039 mg/l 0.010 1 01/04/10 11:15 ND mg/l 0.004 1 01/04/10 11:15 ND mg/l 0.004 1 01/04/10 11:15 ND mg/l 0.010 1 01/04/10 11:15 ND mg/l 0.0002 1 12/30/09 13:20 ND mg/l 0.010 1 01/04/10 11:15 ND mg/l 0.010 1 01/04/10 11:15 ND mg/l 0.007 1 01/04/10 11:15 ND mg/l 0.007 1 01/04/10 11:15 ND mg/l 0.0020 4 12/30/09 09:40 ND mg/l 0.010 1 01/04/10 11:15 | Result Qualifier Units RDL Factor Prepared Analyzed tals - Westborough Lab ND mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:26 ND mg/l 0.005 1 01/04/10 11:15 01/04/10 18:04 0.039 mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.01 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.0002 1 12/30/09 13:20 12/31/09 11:13 ND mg/l 0.0025 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 ND mg/l 0.007 1 01/04/10 11:15 01 | Result Qualifier Units RDL Factor Prepared Analyzed Method tals - Westborough Lab ND mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:26 EPA 3005A ND mg/l 0.005 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.01 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.01 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND mg/l 0.0002 1 12/30/09 13:20 12/31/09 11:13 EPA 7470A ND mg/l 0.025 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A ND </td <td>Result Qualifier Units RDL Factor Prepared Analyzed Method Method tals - Westborough Lab ND mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:26 EPA 3005A 64,6020A ND mg/l 0.005 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B 0.039 mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.01 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.002 1 12/30/09 13:20 12/31/09 11:13 EPA 7470A 64,7470A</td> | Result Qualifier Units RDL Factor Prepared Analyzed Method Method tals - Westborough Lab ND mg/l 0.0020 4 12/30/09 09:40 12/30/09 19:26 EPA 3005A 64,6020A ND mg/l 0.005 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B 0.039 mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.004 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.01 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.010 1 01/04/10 11:15 01/04/10 18:04 EPA 3005A 60,6010B ND mg/l 0.002 1 12/30/09 13:20 12/31/09 11:13 EPA 7470A 64,7470A |



Project Name: WALPOLE PK SOUTH **Lab Number:** L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-04

Client ID: RIZ-3

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 12/28/09 09:03

Date Received: 12/29/09

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:32 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | 0.177 | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:15 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:32 | EPA 3005A | 64,6020A | TD |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 18:07 | EPA 3005A | 60,6010B | Al |



12/28/09 10:00

12/29/09

Project Name: WALPOLE PK SOUTH **Lab Number:** L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L0918777-05

Client ID: RIZ-9

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|------------|-------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | tals - Wes | tborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:38 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:16 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:38 | EPA 3005A | 64,6020A | TD |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 18:10 | EPA 3005A | 60,6010B | AI |



12/28/09 10:43

See Narrative

60,6010B

60,6010B

64,6020A

60,6010B

60,6010B

ΑI

ΑI

TD

ΑI

ΑI

12/29/09

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Date Collected:

Date Received:

01/04/10 11:15 01/04/10 18:14 EPA 3005A

01/04/10 11:15 01/04/10 18:14 EPA 3005A

01/04/10 11:15 01/04/10 18:14 EPA 3005A

Field Prep:

Lab ID: L0918777-06

Client ID: RIZ-10

Sample Location: WALPOLE, MA

ND

ND

ND

ND

ND

Matrix: Water

Selenium, Dissolved

Thallium, Dissolved

Vanadium, Dissolved

Silver, Dissolved

Zinc, Dissolved

| Matrix: | Water | | | | | | | | | |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:44 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.099 | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:55 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:14 | EPA 3005A | 60,6010B | AI |
| | | | | | | | | | | |

1

1

4

1

1

mg/l

mg/l

mg/l

mg/l

mg/l

0.010

0.007

0.0020

0.010

0.050



Project Name: WALPOLE PK SOUTH **Lab Number:** L0918777

Project Number: 12700058 **Report Date:** 01/05/10

SAMPLE RESULTS

Lab ID: L0918777-07

Client ID: MW-3

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 12/28/09 11:50
Date Received: 12/29/09

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-------------|--------------|-------|--------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Me | etals - Wes | stborough La | ab | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:50 | EPA 3005A | 64,6020A | TD |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Barium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 11:56 | EPA 7470A | 64,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | 4 | 12/30/09 09:40 | 12/30/09 19:50 | EPA 3005A | 64,6020A | TD |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 18:17 | EPA 3005A | 60,6010B | Al |



Lab Number:

Project Name: WALPOLE PK SOUTH

L0918777 Project Number: 12700058 **Report Date:** 01/05/10

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|------------------------|----------------------|----------|----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals - | - Westborough Lab fo | r sample | (s): 01- | 07 Batch: | WG395110-1 | | | |
| Antimony, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/09 09:40 | 12/30/09 18:32 | 64,6020A | TD |
| Thallium, Dissolved | ND | mg/l | 0.0005 | 1 | 12/30/09 09:40 | 12/30/09 18:32 | 64,6020A | TD |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|----------------------|----------------------|----------|-----------|--------------------|------------------|------------------|----------------------|----|
| MCP Dissolved Metals | - Westborough Lab fo | r sample | e(s): 01- | 07 Batch: | WG395135-1 | | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 12/30/09 13:20 | 12/31/09 10:53 | 64,7470A | EZ |

Prep Information

Digestion Method: EPA 7470A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------|----------------------|----------|-----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals | - Westborough Lab fo | r sample | (s): 01-0 | 7 Batch: | WG395350-1 | | | |
| Arsenic, Dissolved | ND | mg/l | 0.005 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Barium, Dissolved | ND | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Beryllium, Dissolved | ND | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Cadmium, Dissolved | ND | mg/l | 0.004 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Chromium, Dissolved | ND | mg/l | 0.01 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Lead, Dissolved | ND | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Nickel, Dissolved | ND | mg/l | 0.025 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Selenium, Dissolved | ND | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Silver, Dissolved | ND | mg/l | 0.007 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Vanadium, Dissolved | ND | mg/l | 0.010 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| Zinc, Dissolved | ND | mg/l | 0.050 | 1 | 01/04/10 11:15 | 01/04/10 17:15 | 60,6010B | Al |
| | | | | | | | | |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 **Report Date:** 01/05/10

Method Blank Analysis Batch Quality Control

Prep Information

Digestion Method: EPA 3005A



Project Name: WALPOLE PK SOUTH

Project Number: 12700058

Lab Number: L0918777

| Parameter | | LCS %Recovery | Qua | ı % | LCSD Recover | y Qual | %Recovery Limits | RPD | Qual | RPD Limits | |
|------------------------|-------------------|------------------|---------|-------|-----------------|------------|---------------------|-----|------|------------|--|
| MCP Dissolved Metals - | - Westborough Lab | Associated samp | ole(s): | 01-07 | Batch: | WG395110-2 | WG395110-3 | | | | |
| Antimony, Dissolved | | 97 | | | 94 | | 80-120 | 3 | | 20 | |
| Thallium, Dissolved | | 97 | | | 96 | | 80-120 | 1 | | 20 | |
| MCP Dissolved Metals - | - Westborough Lab | Associated samp | ole(s): | 01-07 | Batch: | WG395135-2 | WG395135-3 | | | | |
| Mercury, Dissolved | | 98 | | | 106 | | 80-120 | 8 | | 20 | |
| MCP Dissolved Metals - | - Westborough Lab | Associated samp | ole(s): | 01-07 | Batch: | WG395350-2 | WG395350-3 | | | | |
| Arsenic, Dissolved | | 110 | | | 109 | | 80-120 | 1 | | 20 | |
| Barium, Dissolved | | 100 | | | 98 | | 80-120 | 2 | | 20 | |
| Beryllium, Dissolved | | 105 | | | 101 | | 80-120 | 4 | | 20 | |
| Cadmium, Dissolved | | 108 | | | 108 | | 80-120 | 0 | | 20 | |
| Chromium, Dissolved | | 100 | | | 105 | | 80-120 | 5 | | 20 | |
| Lead, Dissolved | | 106 | | | 106 | | 80-120 | 0 | | 20 | |
| Nickel, Dissolved | | 101 | | | 101 | | 80-120 | 0 | | 20 | |
| Selenium, Dissolved | | 110 | | | 112 | | 80-120 | 2 | | 20 | |
| Silver, Dissolved | | 103 | | | 105 | | 80-120 | 2 | | 20 | |
| Vanadium, Dissolved | | 102 | | | 104 | | 80-120 | 2 | | 20 | |
| Zinc, Dissolved | | 109 | | | 108 | | 80-120 | 1 | | 20 | |



Project Name: WALPOLE PK SOUTH

Lab Number: L0918777 **Report Date:** 01/05/10 Project Number: 12700058

Sample Receipt and Container Information

YES Were project specific reporting limits specified?

Cooler Information

Cooler **Custody Seal** Α Absent

| Container Info | ormation | | Temp | | | | |
|----------------|----------------------------------|--------|------|-------|------|--------|--|
| Container ID | Container Type | Cooler | рН | deg Ċ | Pres | Seal | Analysis |
| L0918777-01A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-01C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-TL-6020S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-02A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-02B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-02C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-T470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-03A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-03B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-03C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-T470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-04A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-04B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |



Project Name: WALPOLE PK SOUTH Lab Number: L0918777 Project Number: 12700058 **Report Date:** 01/05/10

| Container Info | ormation | Temp | | | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|--|
| Container ID | Container Type | Cooler | рН | deg C | Pres | Seal | Analysis |
| L0918777-04C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-05A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-05B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-05C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-TL-6020S(180),MCP-PB-6010S(180),MCP-PB-6010S(180),MCP-ZN-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-06A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-06B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-06C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-T470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-CN-6010S(180),MCP-CN-6010S(180),MCP-CN-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-07A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-07B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |
| L0918777-07C | Plastic 500ml HNO3 preserved | A | <2 | 5 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SB-6020S(180),MCP-SE-6010S(180),MCP-BE-6010S(180),MCP-NI-6010S(180),MCP-CD-6010S(180),MCP-TL-6020S(180),MCP-TL-6020S(180),MCP-PB-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-AS-6010S(180),MCP-V-6010S(180) |
| L0918777-08A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5 | Υ | Absent | 524.2(14) |

Project Name: WALPOLE PK SOUTH Lab Number: L0918777

Project Number: 12700058 Report Date: 01/05/10

GLOSSARY

Acronyms

EPA · Environmental Protection Agency.

LCS Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD · Laboratory Control Sample Duplicate: Refer to LCS.

MS • Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD · Matrix Spike Sample Duplicate: Refer to MS.

NA · Not Applicable.

NC · Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

ND · Not detected at the reported detection limit for the sample.

NI · Not Ignitable.

RDL - Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A Spectra identified as "Aldol Condensation Product".
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E . Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- ${f P}$ The RPD between the results for the two columns exceeds the method-specified criteria.
- The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RDL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- J : Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).

Report Format: Data Usability Report



Project Name:WALPOLE PK SOUTHLab Number:L0918777Project Number:12700058Report Date:01/05/10

REFERENCES

Methods for the Determination of Organic Compounds in Drinking Water - Supplement II. EPA/600/R-92/129, August 1992.

- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). May 2004.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised December 1, 2009 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. NELAP Accredited Solid Waste/Soil.

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate. Organic Parameters: Haloacetic Acids, Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB).)

Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Calcium Hardness, Silica, Sulfate, Sulfide, Ammonia, Kjeldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH.) Solid Waste/Soil (Inorganic Parameters: Lead in Paint, pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), Reactivity. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH, Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3.3'-Dichlorobenzidine, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9221E, 9222B, 9222D, 9223B, EPA 180.1, 300.0, 353.2, SM2130B, 2320B, 4500Cl-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B,4500NO3-F, EPA 200.7, EPA 200.8, 245.1. Organic Parameters: 504.1, 524.2, SM 6251B.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, Lachat 10-107-06-1-B, SM2320B, 2340B, 2510B, 2540C, 2540D, 426C, 4500Cl-D, 4500Cl-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500P-B.5, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water

Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl)

(EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Fluoride, Sulfate)

353.2 for: Nitrate-N, Nitrite-N; SM4500NO3-F, 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, 2320B, SM2540C. SM4500H-B.

Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics)

(504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), 314.0, 332.

Microbiology Parameters: SM9215B; ENZ. SUB. SM9223; MF-SM9222D

Non-Potable Water

Inorganic Parameters:, (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn)

(EPA 200.7 for: Al,Sb,As,Be,Cd,Cr,Co,Cu,Fe,Pb,Mn,Mo,Ni,Se,Ag,Sr,Tl,Ti,V,Zn,Ca,Mg,Na,K)

245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2540B, 2340B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Ammonia-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-B, C-Titr, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B, E, 5220D, EPA 410.4, SM 5210B,

5310C, 4500CN-CE, 2540D, 4500CL-D, EPA 1664, SM14 510AC, EPA 420.1

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics)

(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCBs-Water), EPA 625 for SVOC Acid Extractables and SVOC Base/Neutral Extractables, 600/4-81-045-PCB-Oil

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM6215B, 9222B, 9223B Colilert, EPA 200.7, 200.8, 245.2, 120.1, 300.0, 314.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 331.0. Organic Parameters: 504.1, 524.2, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 300.0, 350.1, 351.1, 353.2, 420.1, 1664A, SW-846 9010, 9030, 9040B, SM426C, SM2310B, 2540B, 2540D, 4500H+B, 4500NH3-H, 4500NH3-E, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 2320B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-117-07-1-B, LACHAT 10-107-06-1-B, LACHAT 10-107-04-1-J, LACHAT 10-117-07-1-A, SM4500CL-E, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3005A, 3015A, 3510C, 5030B, 8021B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 7.3.3.2, 7.3.4.2, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040, 9045C, 9050C, 1311, 3005A, 3050B, 3051A. Organic Parameters: SW-846 3540C, 3545, 3580A, 5030B, 5035, 8021B, 8260B, 8270C, 8330, 8151A, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 2540C, 2320B, 314.0, SM2120B, 2510B, 5310C, SM4500H-B, EPA 200.8, 245.2. Organic Parameters: 504.1, SM6251B, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.4, SM5220D, 4500Cl-D, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, SM9221CE, 9222D, 9221B, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.1, SM5210B, SW-846 3015, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 3510C, EPA 608, 624, 625, SW-846 5030B, 8021B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 9040B, 3005A, 6010B, 7196A, 5030B, 9010B, 9030B, 1030, 1311, 3050B, 3051, 7471A, 9014, 9012A, 9045C, 9050A, 9065. Organic Parameters: SW-846 8021B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 1311, 1312, 3540C, 3545, 3550B, 3580A, 5035L, 5035H, NJ OQA-QAM-025 Rev.7.)

New York Department of Health Certificate/Lab ID: 11148. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 9215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 314.0, 332.0, SM2320B, EPA 300.0, SM2120B, 4500CN-E, 4500F-C, 4500H-B, 4500NO3-F, 2540C, EPA 120.1, SM 2510B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, 5210B, EPA 410.4, SM5220D, 2310B-4a, 2320B, EPA 200.7, 300.0, LACHAT 10-117-07-1A or B, SM4500Cl-E, 4500F-C, SM15 426C, EPA 350.1, LACHAT 10-107-06-1-B, SM4500NH3-H, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-041-C, SM4500-NO3-F, 4500-NO2-B, 4500P-E, 2540C, 2540B, 2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, S\M3500Cr-D, EPA 245.1, 245.2, 7470A, SM2120B, SM4500-CN-E LACHAT 10-204-00-1-A, EPA 9040B, SM4500-HB, EPA 1664A, SM5310C, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, SM4500S-D, SM5540C, EPA 3005A, 3015. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, EPA 3510C, 5030B, 9010B, 9030B.)

Solid & Hazardous Waste (Inorganic Parameters: 1010, 1030, SW-846 Ch 7 Sec 7.3, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 1312, 3005A, 3050B, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8151A, 8330, 8082, 3540C, 3545, 3546, 3580, 5030B, 5035.)

North Carolina Department of the Environment and Natural Resources <u>Certificate/Lab ID</u>: 666. <u>Organic Parameters</u>: MA-EPH, MA-VPH.

Pennsylvania Department of Environmental Protection Certificate/Lab ID: 68-03671. *NELAP Accredited. Non-Potable Water* (Organic Parameters: EPA 3510C, 5030B, 625, 624. 608, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1010, 1030, 1311, 3050B, 3051, 6010B, EPA 7.3.3.2, EPA 7.3.4.2, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065. Organic Parameters: 3540C, 3545, 3580A, 5035, 8021B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Rhode Island Department of Health Certificate/Lab ID: LAO00065. *NELAP Accredited via NY-DOH.*Refer to MA-DEP Certificate for Potable and Non-Potable Water.
Refer to NY-DOH Certificate for Potable and Non-Potable Water.

Texas Commisson on Environmental Quality Certificate/Lab ID: T104704476-09-1. **NELAP Accredited.** Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664, 200.7, 200.8, 245.1, 245.2, 300.0, 350.1, 351.1, 353.2, 376.2, 410.4, 420.1, 6010, 6020, 7196, 7470, 9040, SM 2120B, 2310B, 2320B, 2510B, 2540B, 2540B, 2540D, 426C, 4500CL-E, 4500CN-E, 4500F-C, 4500H+B, 4500NH3-H, 4500NO2B, 4500P-E, 4500 S2⁻ D, 510C, 5210B, 5220D, 5310C, 5540C. Organic Parameters: EPA 608, 624, 625, 8081, 8082, 8151, 8260, 8270, 8330.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 9012, 9014, 9040, 9045, 9050, 9065.)

Utah Department of Health Certificate/Lab ID: AAMA. **NELAP Accredited.** Non-Potable Water (Inorganic Parameters: Chloride EPA 300.0)

Department of Defense Certificate/Lab ID: L2217.

Drinking Water (Inorganic Parameters: SM 4500H-B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: EPA 200.7, 200.8, 6010B, 6020, 245.1, 245.2, 7470A, 9040B, 300.0, 9251, 9038, 350.1, 353.2, 351.1, 314, 120.1, 9050A, 410.4, 9060, 1664, 420.1, LACHAT 10-107-06-1-B, SM 4500CN-E, 4500H-B, 4500CL-E, 4500F-BC, 4500SO4-E, 426C, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500NO2-B, 4500Norg-C, 4500PE, 2510B, 5540C, 5220D, 5310C, 2540B, 2540C, 2540D, 510C, 4500S2-AD, 3005A, 3015, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8330, 625, 8082, 8151A, 8081A, 3510C, 5030B.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 200.7, 6010B, 7471A, 9040B, 9045C, 9065, 420.1, 9012A, 6860, 1311, 1312, 3050B, 9030B, 3051, 9010B, 3540C, SM 510ABC, 4500CN-CE, 2540G, SW-846 7.3, Organic Parameters: EPA 8260B, 8270C, 8330, 8082, 8081A, 8151A, 3545, 3546, 3580, 5035.)

Analytes Not Accredited by NELAP

Certification is not available by NELAP for the following analytes: **EPA 8260B**: Freon-113, 1,2,4,5-Tetramethylbenzene. **EPA 8330A**: PETN, Picric Acid, Nitroglycerine, 2,6-DANT, 2,4-DANT. **EPA 8270C**: Methyl naphthalene, Dimethyl naphthalene, Total Methylnapthalenes, Total Dimethylnaphthalenes, 1,4-Diphenylhydrazine (Azobenzene). **EPA 625**: 4-Chloroaniline. **EPA 350.1** for Ammonia in a Soil matrix.

| IS YOUR PROJECT MA MCP or CT RCP? FORM NO: 01-01 (rev. 14-0CT-07) | 8 Tripbluc | 6 RIZ-10 | 9-7H7 E | 18777.1 Mw-9 | ALPHA Lab ID (Lab Use Only) Sample ID | Other Project specific Requirements/Comments/Detection Limits $DL/ \neq \mathit{KCEO}-1$ | Email: IAM. CAMMEN Athorped. Com These samples have been previously analyzed by Alpha | Fax: 508 903 2001 | Fanu | Address: D. 1 1 CL | 508-898-9193 Int Information | CHAIN (AMAZICA) TEL: 508-832-9300 TEL: 508-832-9300 |
|--|---------------------|----------|------------------------|---------------------------------------|--|--|--|--|--|--|---------------------------------|---|
| Relinquished By: Respression 12/16/9 - 8/0K Apha | 643 | 10/28/4 0725 | 12/21/9 1012 6W 180 | Collection Sample Sampler's Date Time Matrix Initials | | pha Date Due: \\(\(\) \\(\) \\\(\) \\\(\) \\\\\(\) \\\\\(\) \\\\\(\) \\\\\(\) \\\\\\\\ | Turn-Around Time | 1 2000 | Project # 12 7005 8 Project Manager: Ra. Ch. In Toure | Project Name: Unipole IVE SOUL | PAGE |
| HLI N Rederved By: 12/24/C | × < | | | × × × × × × × × × × × × × × × × × × × | ers Vo | ANA C. SZY MCD 14 M | etak | XYes ☐ No Are MCP Analytical Methods Required? ☐ Yes | MA MCP PR | Regulatory Requirements/Report Limits State /Fed Program Criteria | | ed in Lab: (ヨ)ユ의 0 句 Information - Data Deliverables |
| Please print clearly, legibly and completely. Samples cannot be logged in and turnaround time clock will not start until any ambiguities are resolved All samples submitted are subject to the phase Terms and Conditions of See reverse side | ~ * | | vsix o. vs.un Pitter 2 | metals fell-filled 3 | (Please specify below) Sample Specific Comments | eded do | SAMPLE HANDLING Filtration | quired? dence Protocols) Required? | ESUMPTIVE CERTAINTY - CT REASONABLE CONFIDENCE PROTOCOLS | | a same as Client into PO#: | |



ANALYTICAL REPORT

Lab Number: L1004931

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: Ian Cannan Phone: (508) 903-2039

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Report Date: 04/15/10

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Number: 12700058 **Report Date:** 04/15/10

| Alpha Sample ID | Client ID | Sample Location | Collection Date/Time |
|--------------------|-----------|--------------------|-------------------------|
| L1004931-01 | GHC-5 | WALPOLE, MA | 04/07/10 12:32 |
| L1004931-02 | GHC-5-A | WALPOLE, MA | 04/07/10 12:56 |
| L1004931-03 | GHC-5-B | WALPOLE, MA | 04/07/10 12:57 |
| L1004931-04 | GHC-5-C | WALPOLE, MA | 04/07/10 12:58 |
| L1004931-05 | RIZ-2 | WALPOLE, MA | 04/07/10 13:39 |
| L1004931-06 | RIZ-2-A | WALPOLE, MA | 04/07/10 13:55 |
| L1004931-07 | RIZ-2-B | WALPOLE, MA | 04/07/10 13:56 |
| L1004931-08 | RIZ-2-C | WALPOLE, MA | 04/07/10 13:57 |
| L1004931-09 | RIZ-10 | WALPOLE, MA | 04/07/10 14:19 |
| L1004931-10 | RIZ-10-A | WALPOLE, MA | 04/07/10 14:32 |
| L1004931-11 | RIZ-10-B | WALPOLE, MA | 04/07/10 14:33 |
| L1004931-12 | RIZ-10-C | WALPOLE, MA | 04/07/10 14:34 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP Analytical Methods.

| Α | Were all samples received by the laboratory in a condition consistent with those described on their Chain-of-Custody documentation for the data set? | YES |
|------|---|-----|
| В | Were all QA/QC procedures required for the specified analytical methods(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | YES |
| С | Does the analytical data included in this report meet all the requirements for "Presumptive Certainty", as described in section 2.0 of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"? | YES |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications, as specified in Section 11.3? | N/A |
| A re | sponse to questions E and F is required for "Presumptive Certainty" status | |
| E | Were all QC performance standards and recommendations for the specified method(s) achieved? | YES |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | NO |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



L1004931

Lab Number:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 **Report Date:** 04/15/10

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

MCP Related Narratives

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

For additional information, please contact Client Services at 800-624-9220.

The date collected for L1004931-01 through -04 was obtained from the sample labels.

Metals

In reference to question F:

All samples were analyzed for a subset of MCP elements per the Chain of Custody.



Project Number: 12700058 **Report Date:** 04/15/10

Case Narrative (continued)

Non-MCP Related Narratives

Phenolics, Total

L1004931-01 and -09 have elevated detection limits due to the dilutions required by the sample matrices.

Coliform, Fecal (MF)

L1004931-02, -03, -04, -10, -11 and -12 have elevated detection limits due to the dilutions required by the sample matrices.

L1004931-06, -07 and -08 have elevated detection limits due to the dilutions required by the method.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative

ALPHA

Date: 04/15/10

ORGANICS



VOLATILES



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: Report Date: 12700058 04/15/10

SAMPLE RESULTS

Lab ID: Date Collected: L1004931-01 04/07/10 12:32

Client ID: GHC-5

Date Received: 04/07/10 Field Prep: Sample Location: WALPOLE, MA See Narrative

Matrix: Water Analytical Method: 16,524.2 04/09/10 17:33 Analytical Date:

Analyst: TT

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | gh Lab | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| Chloroform | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Bromoform | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Benzene | ND | | ug/l | 0.50 | 1 |
| Toluene | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Chloromethane | ND | | ug/l | 0.50 | 1 |
| Bromomethane | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| Chloroethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: Report Date: 12700058 04/15/10

SAMPLE RESULTS

Lab ID: Date Collected: L1004931-01 04/07/10 12:32

Client ID: GHC-5 Date Received: 04/07/10

Sample Location: WALPOLE, MA Field Prep: See Narrative Parameter RDL **Dilution Factor** Qualifier Units Result

| Parameter | Result | Qualifier Units | KDL | Dilution Factor |
|-------------------------------------|------------|-----------------|------|-----------------|
| Volatile Organics by GC/MS - Westbo | orough Lab | | | |
| 1,3-Dichlorobenzene | ND | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | ug/l | 0.50 | 1 |
| Styrene | ND | ug/l | 0.50 | 1 |
| o-Xylene | ND | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | ug/l | 0.50 | 1 |
| Naphthalene | ND | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | ug/l | 0.50 | 1 |
| Bromobenzene | ND | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | ug/l | 0.50 | 1 |
| Dibromomethane | ND | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | ug/l | 0.50 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 | |
| 4-Bromofluorobenzene | 89 | | 80-120 | |



04/07/10

See Narrative

Date Received:

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-05 Date Collected: 04/07/10 13:39

Client ID: RIZ-2

Sample Location: WALPOLE, MA Field Prep:

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 04/09/10 18:05

Analyst: TT

| Volatile Organics by GC/MS - Westborough Lab Methylene chloride ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichklorofucorethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane | Volatile Organics by GC/MS - Westborou | gh Lab | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itaris-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bernache ND ug/l 0.50 1 Bernacher ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 </td <td>1,1-Dichloroethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromodichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromofichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Eenzene ND ug/l 0.50 1 Tolluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorotluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Chloro | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itrans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <td< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| ND | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ehylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND ug/l 0.50 1 | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-05 Date Collected: 04/07/10 13:39

Client ID: RIZ-2 Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL **Dilution Factor** Parameter Qualifier Units Result Volatile Organics by GC/MS - Westborough Lab ND 1,3-Dichlorobenzene ug/l 0.50 1 ND 0.50 1 1,4-Dichlorobenzene ug/l Styrene ND 0.50 ug/l 1 o-Xylene ND ug/l 0.50 1 ND 0.50 1,1-Dichloropropene ug/l 1 ND 0.50 2,2-Dichloropropane ug/l 1 1,1,1,2-Tetrachloroethane ND 0.50 1 ug/l ND 0.50 1 1,2,3-Trichloropropane ug/l Bromochloromethane ND ug/l 0.50 1 ND 0.50 n-Butylbenzene ug/l 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND 0.50 ug/l 1 tert-Butylbenzene ND ug/l 0.50 1 ND 0.50 1 1,2,3-Trichlorobenzene ug/l ND 0.50 1 1,2,4-Trichlorobenzene ug/l 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND 0.50 1 ug/l Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND 0.50 1 ug/l ND p-Chlorotoluene ug/l 0.50 1

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 104 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 80-120 | |

ND

ND

ND

ND

ND

ug/l

ug/l

ug/l

ug/l

ug/l

0.50

0.50

0.50

0.50

0.50



1

1

1

1

1

Dibromomethane

1,2-Dibromoethane

1,3-Dichloropropane

Methyl tert butyl ether

1,2-Dibromo-3-chloropropane

04/07/10

See Narrative

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09 Date Collected: 04/07/10 14:19

Client ID: RIZ-10

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 04/09/10 18:38

Analyst: TT

| Volatile Organics by GC/MS - Westborough Lab Methylene chloride ND ug/l 0.50 1 1,1-Dichloroethane ND ug/l 0.50 1 Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichklorofucorethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromodichkoromethane ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform | Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|--|--------|-----------|-------|------|-----------------|
| 1,1-Dichloroethane | Volatile Organics by GC/MS - Westborou | gh Lab | | | | |
| Chloroform ND ug/l 0.50 1 Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | Methylene chloride | ND | | ug/l | 0.50 | 1 |
| Carbon tetrachloride ND ug/l 0.50 1 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itaris-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bernache ND ug/l 0.50 1 Bernacher ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 </td <td>1,1-Dichloroethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | 1,1-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloropropane ND ug/l 0.50 1 Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 Trichloroethane ND ug/l 0.50 1 1,1-1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromodichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | Chloroform | ND | | ug/l | 0.50 | 1 |
| Dibromochloromethane ND ug/l 0.50 1 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Bromofichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 <td>Carbon tetrachloride</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>1</td> | Carbon tetrachloride | ND | | ug/l | 0.50 | 1 |
| 1,1,2-Trichloroethane ND ug/l 0.50 1 Tetrachloroethane ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichloroffuoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Eenzene ND ug/l 0.50 1 Tolluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 | 1,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Tetrachloroethene ND ug/l 0.50 1 Chlorobenzene ND ug/l 0.50 1 Trichlorotluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Chloro | Dibromochloromethane | ND | | ug/l | 0.50 | 1 |
| Chlorobenzene ND ug/l 0.50 1 Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 P/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| Trichlorofluoromethane ND ug/l 0.50 1 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 Itrans-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 | Tetrachloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichloroethane ND ug/l 0.50 1 1,1,1-Trichloroethane ND ug/l 0.50 1 Bromodichloromethane ND ug/l 0.50 1 trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 Ethyloenzene ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 <td< td=""><td>Chlorobenzene</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>1</td></td<> | Chlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,1,1-Trichloroethane | Trichlorofluoromethane | ND | | ug/l | 0.50 | 1 |
| ND | 1,2-Dichloroethane | ND | | ug/l | 0.50 | 1 |
| trans-1,3-Dichloropropene ND ug/l 0.50 1 cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 | 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,3-Dichloropropene ND ug/l 0.50 1 Bromoform ND ug/l 0.50 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 Ehylbenzene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromodichloromethane | ND | | ug/l | 0.50 | 1 |
| Bromoform ND ug/l 0.50 1 | trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 1,1,2,2-Tetrachloroethane ND ug/l 0.50 1 Benzene ND ug/l 0.50 1 Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| Benzene ND ug/l 0.50 1 | Bromoform | ND | | ug/l | 0.50 | 1 |
| Toluene ND ug/l 0.50 1 Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| Ethylbenzene ND ug/l 0.50 1 p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Benzene | ND | | ug/l | 0.50 | 1 |
| p/m-Xylene ND ug/l 0.50 1 Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Toluene | ND | | ug/l | 0.50 | 1 |
| Chloromethane ND ug/l 0.50 1 Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Ethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromomethane ND ug/l 0.50 1 Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | p/m-Xylene | ND | | ug/l | 0.50 | 1 |
| Vinyl chloride ND ug/l 0.50 1 Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloromethane | ND | | ug/l | 0.50 | 1 |
| Chloroethane ND ug/l 0.50 1 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Bromomethane | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloroethene ND ug/l 0.50 1 trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Vinyl chloride | ND | | ug/l | 0.50 | 1 |
| trans-1,2-Dichloroethene ND ug/l 0.50 1 cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | Chloroethane | ND | | ug/l | 0.50 | 1 |
| cis-1,2-Dichloroethene ND ug/l 0.50 1 Trichloroethene ND ug/l 0.50 1 | 1,1-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| Trichloroethene ND ug/l 0.50 1 | trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| | cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | 1 |
| 1,2-Dichlorobenzene ND ug/l 0.50 1 | Trichloroethene | ND | | ug/l | 0.50 | 1 |
| | 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09 Date Collected: 04/07/10 14:19

Client ID: RIZ-10 Date Received: 04/07/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|--|--------|-----------|-------|------|-----------------|
| Volatile Organics by GC/MS - Westborough | n Lab | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | 1 |
| Styrene | ND | | ug/l | 0.50 | 1 |
| o-Xylene | ND | | ug/l | 0.50 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | 1 |
| Naphthalene | ND | | ug/l | 0.50 | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 104 | | 80-120 | |
| 4-Bromofluorobenzene | 89 | | 80-120 | |



Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 04/09/10 13:54

Analyst: TT

| arameter | Result | Qualifier | Units | RDL | |
|-----------------------------|----------------|------------------|----------|--------|------------|
| platile Organics by GC/MS - | Westborough La | b for sample(s): | 01,05,09 | Batch: | WG407461-2 |
| Methylene chloride | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | |
| Chloroform | ND | | ug/l | 0.50 | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | |
| Dibromochloromethane | ND | | ug/l | 0.50 | |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | |
| Tetrachloroethene | ND | | ug/l | 0.50 | |
| Chlorobenzene | ND | | ug/l | 0.50 | |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | |
| Bromodichloromethane | ND | | ug/l | 0.50 | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| Bromoform | ND | | ug/l | 0.50 | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | |
| Benzene | ND | | ug/l | 0.50 | |
| Toluene | ND | | ug/l | 0.50 | |
| Ethylbenzene | ND | | ug/l | 0.50 | |
| p/m-Xylene | ND | | ug/l | 0.50 | |
| Chloromethane | ND | | ug/l | 0.50 | |
| Bromomethane | ND | | ug/l | 0.50 | |
| Vinyl chloride | ND | | ug/l | 0.50 | |
| Chloroethane | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| Trichloroethene | ND | | ug/l | 0.50 | |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | |



Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 04/09/10 13:54

Analyst: TT

| arameter | Result | Qualifier | Units | RDL | |
|-------------------------------|----------------|------------------|------------|--------|------------|
| olatile Organics by GC/MS - V | Vestborough La | b for sample(s): | : 01,05,09 | Batch: | WG407461-2 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | |
| Styrene | ND | | ug/l | 0.50 | |
| o-Xylene | ND | | ug/l | 0.50 | |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | |
| Bromochloromethane | ND | | ug/l | 0.50 | |
| n-Butylbenzene | ND | | ug/l | 0.50 | |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | |
| Isopropylbenzene | ND | | ug/l | 0.50 | |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | |
| Naphthalene | ND | | ug/l | 0.50 | |
| n-Propylbenzene | ND | | ug/l | 0.50 | |
| sec-Butylbenzene | ND | | ug/l | 0.50 | |
| tert-Butylbenzene | ND | | ug/l | 0.50 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | |
| Bromobenzene | ND | | ug/l | 0.50 | |
| o-Chlorotoluene | ND | | ug/l | 0.50 | |
| p-Chlorotoluene | ND | | ug/l | 0.50 | |
| Dibromomethane | ND | | ug/l | 0.50 | |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | |
| | | | | | |



Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 04/09/10 13:54

Analyst: TT

Parameter Result Qualifier Units RDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01,05,09 Batch: WG407461-2

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | Acceptance | | | | | | | | |
|------------------------|------------|-----------|----------|--|--|--|--|--|--|
| Surrogate | %Recovery | Qualifier | Criteria | | | | | | |
| | | | | | | | | | |
| 1,2-Dichlorobenzene-d4 | 102 | | 80-120 | | | | | | |
| 4-Bromofluorobenzene | 93 | | 80-120 | | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| arameter | LCS %Recovery Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | Qual | RPD Limits |
|---|-----------------------|---------------------|--------------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough I | ab Associated sample | (s): 01,05,09 Batch | : WG407461-1 | | | |
| Methylene chloride | 97 | - | 70-130 | - | | |
| 1,1-Dichloroethane | 101 | - | 70-130 | - | | |
| Chloroform | 104 | - | 70-130 | - | | |
| Carbon tetrachloride | 98 | - | 70-130 | - | | |
| 1,2-Dichloropropane | 102 | - | 70-130 | - | | |
| Dibromochloromethane | 113 | - | 70-130 | - | | |
| 1,1,2-Trichloroethane | 120 | - | 70-130 | - | | |
| Tetrachloroethene | 119 | - | 70-130 | - | | |
| Chlorobenzene | 94 | - | 70-130 | - | | |
| Trichlorofluoromethane | 97 | - | 70-130 | - | | |
| 1,2-Dichloroethane | 102 | - | 70-130 | - | | |
| 1,1,1-Trichloroethane | 102 | - | 70-130 | - | | |
| Bromodichloromethane | 98 | - | 70-130 | - | | |
| trans-1,3-Dichloropropene | 115 | - | 70-130 | - | | |
| cis-1,3-Dichloropropene | 111 | - | 70-130 | - | | |
| Bromoform | 88 | - | 70-130 | - | | |
| 1,1,2,2-Tetrachloroethane | 95 | - | 70-130 | - | | |
| Benzene | 106 | - | 70-130 | - | | |
| Toluene | 125 | - | 70-130 | - | | |
| Ethylbenzene | 92 | - | 70-130 | - | | |
| p/m-Xylene | 94 | - | 70-130 | - | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| Parameter | LCS %Recovery | Qual | LCSI %Recov | | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|------------|----------------|-------|---------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough L | ab Associated | sample(s): | 01,05,09 | Batch | : WG407 | 7461-1 | | | |
| Chloromethane | 96 | | - | | | 70-130 | - | | |
| Bromomethane | 93 | | - | | | 70-130 | - | | |
| Vinyl chloride | 106 | | - | | | 70-130 | - | | |
| Chloroethane | 88 | | - | | | 70-130 | - | | |
| 1,1-Dichloroethene | 96 | | - | | | 70-130 | - | | |
| trans-1,2-Dichloroethene | 97 | | - | | | 70-130 | - | | |
| cis-1,2-Dichloroethene | 107 | | - | | | 70-130 | - | | |
| Trichloroethene | 101 | | - | | | 70-130 | - | | |
| 1,2-Dichlorobenzene | 105 | | - | | | 70-130 | - | | |
| 1,3-Dichlorobenzene | 104 | | - | | | 70-130 | - | | |
| 1,4-Dichlorobenzene | 104 | | - | | | 70-130 | - | | |
| Styrene | 94 | | - | | | 70-130 | - | | |
| o-Xylene | 94 | | - | | | 70-130 | - | | |
| 1,1-Dichloropropene | 103 | | - | | | 70-130 | - | | |
| 2,2-Dichloropropane | 102 | | - | | | 70-130 | - | | |
| 1,1,1,2-Tetrachloroethane | 88 | | - | | | 70-130 | - | | |
| 1,2,3-Trichloropropane | 84 | | - | | | 70-130 | - | | |
| Bromochloromethane | 109 | | - | | | 70-130 | - | | |
| n-Butylbenzene | 104 | | - | | | 70-130 | - | | |
| Dichlorodifluoromethane | 90 | | - | | | 70-130 | - | | |
| Hexachlorobutadiene | 103 | | - | | | 70-130 | - | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| arameter | LCS %Recovery | Qual | LCSD %Recovery | / Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|-----------|-------------------|-----------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough L | ab Associated s | ample(s): | 01,05,09 Ba | atch: WG4 | 107461-1 | | | |
| Isopropylbenzene | 92 | | - | | 70-130 | - | | |
| p-isopropyltoluene | 91 | | - | | 70-130 | - | | |
| Naphthalene | 93 | | - | | 70-130 | - | | |
| n-Propylbenzene | 91 | | - | | 70-130 | - | | |
| sec-Butylbenzene | 92 | | - | | 70-130 | - | | |
| tert-Butylbenzene | 92 | | - | | 70-130 | - | | |
| 1,2,3-Trichlorobenzene | 102 | | - | | 70-130 | - | | |
| 1,2,4-Trichlorobenzene | 101 | | - | | 70-130 | - | | |
| 1,2,4-Trimethylbenzene | 94 | | - | | 70-130 | - | | |
| 1,3,5-Trimethylbenzene | 92 | | - | | 70-130 | - | | |
| Bromobenzene | 92 | | - | | 70-130 | - | | |
| o-Chlorotoluene | 91 | | - | | 70-130 | - | | |
| p-Chlorotoluene | 93 | | - | | 70-130 | - | | |
| Dibromomethane | 105 | | - | | 70-130 | - | | |
| 1,2-Dibromoethane | 114 | | - | | 70-130 | - | | |
| 1,2-Dibromo-3-chloropropane | 99 | | - | | 70-130 | - | | |
| 1,3-Dichloropropane | 128 | | - | | 70-130 | - | | |
| Methyl tert butyl ether | 95 | | - | | 70-130 | - | | |

Lab Number: L1004931

Report Date: 04/15/10

| | LCS | | LCSD | | %Recovery | | | |
|-----------|-----------|------|-----------|------|-----------|-----|------|------------|
| Parameter | %Recovery | Qual | %Recovery | Qual | Limits | RPD | Qual | RPD Limits |
| | | | | | | | | |

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01,05,09 Batch: WG407461-1

WALPOLE PARK SOUTH

12700058

| Surrogate | LCS %Recovery | Qual | LCSD %Recovery | Qual | Acceptance Criteria |
|------------------------|------------------|------|-------------------|------|------------------------|
| 1,2-Dichlorobenzene-d4 | 98 | | | | 80-120 |
| 4-Bromofluorobenzene | 89 | | | | 80-120 |



Project Name:

Project Number:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | | Recovery Limits | RPD | RPD Qual Limits |
|-----------------------------|------------------|-------------|---------------|-----------------|--------|--------------|------------------|---------------------|--------------------|--------|--------------------|
| olatile Organics by GC/MS - | - Westborough | n Lab Assoc | ciated sample | e(s): 01,05,09 | QC Bat | tch ID: WG | 3407461-5 (| QC Sam _l | ple: L1004 | 819-01 | Client ID: MS |
| Methylene chloride | ND | 4 | 3.9 | 97 | | - | - | | 70-130 | - | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.4 | 111 | | - | - | | 70-130 | - | 20 |
| Chloroform | ND | 4 | 4.7 | 117 | | - | - | | 70-130 | - | 20 |
| Carbon tetrachloride | ND | 4 | 4.8 | 120 | | - | - | | 70-130 | - | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.6 | 114 | | - | - | | 70-130 | - | 20 |
| Dibromochloromethane | ND | 4 | 4.9 | 123 | | - | - | | 70-130 | - | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 5.4 | 136 | Q | - | - | | 70-130 | - | 20 |
| Tetrachloroethene | ND | 4 | 5.4 | 136 | Q | - | - | | 70-130 | - | 20 |
| Chlorobenzene | ND | 4 | 4.6 | 115 | | - | - | | 70-130 | - | 20 |
| Trichlorofluoromethane | ND | 4 | 4.6 | 115 | | - | - | | 70-130 | - | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.6 | 115 | | - | - | | 70-130 | - | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 4.9 | 123 | | - | - | | 70-130 | - | 20 |
| Bromodichloromethane | ND | 4 | 4.4 | 109 | | - | - | | 70-130 | - | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 4.4 | 111 | | - | - | | 70-130 | - | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 5.0 | 124 | | - | - | | 70-130 | - | 20 |
| Bromoform | ND | 4 | 4.0 | 100 | | - | - | | 70-130 | - | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 4.6 | 114 | | - | - | | 70-130 | - | 20 |
| Benzene | ND | 4 | 4.4 | 111 | | - | - | | 70-130 | - | 20 |
| Toluene | ND | 4 | 5.3 | 133 | Q | - | - | | 70-130 | - | 20 |
| Ethylbenzene | ND | 4 | 4.4 | 111 | | - | - | | 70-130 | - | 20 |
| p/m-Xylene | ND | 8 | 8.9 | 111 | | - | - | | 70-130 | - | 20 |
| | | | | | | | | | | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | | Recovery Limits | RPD | RPD Qual Limits |
|--------------------------------------|------------------|-------------|--------------|-----------------|--------|--------------|------------------|---------|--------------------|--------|--------------------|
| /olatile Organics by GC/MS Sample | - Westborough | n Lab Assoc | iated sample | (s): 01,05,09 | QC Bat | ch ID: WG | 3407461-5 C | QC Samp | ole: L1004 | 819-01 | Client ID: MS |
| Chloromethane | ND | 4 | 5.2 | 131 | Q | - | - | | 70-130 | - | 20 |
| Bromomethane | ND | 4 | 4.0 | 101 | | - | - | | 70-130 | - | 20 |
| Vinyl chloride | ND | 4 | 6.9 | 173 | Q | - | - | | 70-130 | - | 20 |
| Chloroethane | ND | 4 | 4.3 | 107 | | - | - | | 70-130 | - | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.1 | 103 | | - | - | | 70-130 | - | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.2 | 105 | | - | - | | 70-130 | - | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 4.7 | 118 | | - | - | | 70-130 | - | 20 |
| Trichloroethene | ND | 4 | 4.4 | 110 | | - | - | | 70-130 | - | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 4.3 | 108 | | - | - | | 70-130 | - | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 4.2 | 106 | | - | - | | 70-130 | - | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 4.1 | 103 | | - | - | | 70-130 | - | 20 |
| Styrene | ND | 4 | 4.2 | 105 | | - | - | | 70-130 | - | 20 |
| o-Xylene | ND | 4 | 4.4 | 111 | | - | - | | 70-130 | - | 20 |
| 1,1-Dichloropropene | ND | 4 | 4.7 | 118 | | - | - | | 70-130 | - | 20 |
| 2,2-Dichloropropane | ND | 4 | 4.7 | 119 | | - | - | | 70-130 | - | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 4.0 | 101 | | - | - | | 70-130 | - | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 4.2 | 106 | | - | - | | 70-130 | - | 20 |
| Bromochloromethane | ND | 4 | 4.6 | 115 | | - | - | | 70-130 | - | 20 |
| n-Butylbenzene | ND | 4 | 4.5 | 113 | | - | - | | 70-130 | - | 20 |
| Dichlorodifluoromethane | ND | 4 | 5.6 | 140 | Q | - | - | | 70-130 | - | 20 |
| Hexachlorobutadiene | ND | 4 | 3.9 | 99 | | - | - | | 70-130 | - | 20 |
| | | | | | | | | | | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | | Recovery Limits | RPD | Qual | RPD Limits |
|----------------------------------|------------------|-------------|-------------|-----------------|----------|--------------|------------------|--------|--------------------|--------|--------|---------------|
| olatile Organics by GC/MS cample | - Westborough | Lab Associ | ated sample | (s): 01,05,09 | QC Bato | ch ID: WG | 407461-5 (| QC Sam | ple: L1004 | 819-01 | Client | ID: MS |
| Isopropylbenzene | ND | 4 | 3.9 | 97 | | - | - | | 70-130 | - | | 20 |
| p-Isopropyltoluene | ND | 4 | 4.2 | 104 | | - | - | | 70-130 | - | | 20 |
| Naphthalene | ND | 4 | 3.8 | 95 | | - | - | | 70-130 | - | | 20 |
| n-Propylbenzene | ND | 4 | 4.5 | 112 | | - | - | | 70-130 | - | | 20 |
| sec-Butylbenzene | ND | 4 | 4.4 | 110 | | - | - | | 70-130 | - | | 20 |
| tert-Butylbenzene | ND | 4 | 4.2 | 106 | | - | - | | 70-130 | - | | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.9 | 98 | | - | - | | 70-130 | - | | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.6 | 91 | | - | - | | 70-130 | - | | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 4.4 | 110 | | - | - | | 70-130 | - | | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 4.3 | 108 | | - | - | | 70-130 | - | | 20 |
| Bromobenzene | ND | 4 | 4.3 | 107 | | - | - | | 70-130 | - | | 20 |
| o-Chlorotoluene | ND | 4 | 4.3 | 107 | | - | - | | 70-130 | - | | 20 |
| p-Chlorotoluene | ND | 4 | 4.2 | 106 | | - | - | | 70-130 | - | | 20 |
| Dibromomethane | ND | 4 | 4.3 | 109 | | - | - | | 70-130 | - | | 20 |
| 1,2-Dibromoethane | ND | 4 | 4.2 | 105 | | - | - | | 70-130 | - | | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 4.0 | 101 | | - | - | | 70-130 | - | | 20 |
| 1,3-Dichloropropane | ND | 4 | 5.3 | 134 | Q | - | - | | 70-130 | - | | 20 |
| Methyl tert butyl ether | ND | 4 | 4.1 | 103 | <u> </u> | - | - | | 70-130 | - | | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L1004931

Report Date:

04/15/10

| | Native | MS | MS | MS | | MSD | MSD | | Recovery | | | RPD |
|-----------|--------|-------|-------|-----------|------|-------|-----------|------|----------|-----|------|--------|
| Parameter | Sample | Added | Found | %Recovery | Qual | Found | %Recovery | Qual | Limits | RPD | Qual | Limits |

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01,05,09 QC Batch ID: WG407461-5 QC Sample: L1004819-01 Client ID: MS Sample

| | MS | MSD | Acceptance | |
|------------------------|----------------------|----------------------|------------|--|
| Surrogate | % Recovery Qualifier | % Recovery Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 99 | | 80-120 | |
| 4-Bromofluorobenzene | 100 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L1004931

Report Date:

04/15/10

| arameter | Native Sample | Duplicate Sample | Units | RPD | Qual RPD Limits |
|--|--------------------------------|-------------------|------------|--------------|-------------------------|
| olatile Organics by GC/MS - Westborough Lab ample | Associated sample(s): 01,05,09 | QC Batch ID: WG40 |)7461-6 QC | Sample: L100 | 04819-02 Client ID: DUP |
| Methylene chloride | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethane | ND | ND | ug/l | NC | 20 |
| Chloroform | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene | ND | ND | ug/l | NC | 20 |
| Chlorobenzene | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| Bromoform | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| Benzene | ND | ND | ug/l | NC | 20 |
| Toluene | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01,05,09 Bample QC Batch ID: WG407461-6 QC Sample: L1004819-02 Client ID: DUSAmple Ethylbenzere ND ND ND ug1 NC 20 p/m-Xylene ND ND ND ug1 NC 20 Chloromethane ND ND ND ug1 NC 20 Bromomethane ND ND ND ug1 NC 20 Vinyl chloride ND ND ND ug1 NC 20 Chloroethane ND ND ND ug1 NC 20 1,1-Dichloroethene ND ND ND ug1 NC 20 trans-1,2-Dichloroethene ND ND ND ug1 NC 20 Trichloroethene ND ND ND ug1 NC 20 1,2-Dichloroethezene ND ND ND ug1 NC 20 1,3-Dichloroethezene ND <t< th=""><th>Parameter</th><th>Native Sample</th><th>Duplicate Sample</th><th>Units</th><th>RPD</th><th>RPD Limits</th></t<> | Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|---------------------------|--------------------------------|-------------------|------------|----------------|----------------------|
| p/m-Xylene ND ND ug/l NC 20 Chloromethane ND ND ND ug/l NC 20 Bromomethane ND ND ND ug/l NC 20 Vinyl chloride ND ND ND ug/l NC 20 Chloroethane ND ND ND ug/l NC 20 1,1-Dichloroethane ND ND ND ug/l NC 20 trans-1,2-Dichloroethane ND ND ND ug/l NC 20 trans-1,2-Dichloroethane ND ND ND ug/l NC 20 Trichloroethane ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC | | Associated sample(s): 01,05,09 | QC Batch ID: WG40 | 07461-6 QC | Sample: L10048 | 19-02 Client ID: DUP |
| Chloromethane ND ND ug/l NC 20 Bromomethane ND ND ND ug/l NC 20 Vinyl chloride ND ND ND ug/l NC 20 Chloroethane ND ND ND ug/l NC 20 1,1-Dichloroethene ND ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ND ug/l NC 20 dis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l | Ethylbenzene | ND | ND | ug/l | NC | 20 |
| Bromomethane ND ND ug/l NC 20 Vinyl chloride ND ND ND ug/l NC 20 Chloroethane ND ND ND ug/l NC 20 1,1-Dichloroethene ND ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l | p/m-Xylene | ND | ND | ug/l | NC | 20 |
| Vinyl chloride ND ND ug/l NC 20 Chloroethane ND ND ND ug/l NC 20 1,1-Dichloroethene ND ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 0-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l <t< td=""><td>Chloromethane</td><td>ND</td><td>ND</td><td>ug/l</td><td>NC</td><td>20</td></t<> | Chloromethane | ND | ND | ug/l | NC | 20 |
| Chloroethane ND ND ug/l NC 20 1,1-Dichloroethene ND ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 0-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l | Bromomethane | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloroethene ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 0-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ND | Vinyl chloride | ND | ND | ug/l | NC | 20 |
| trans-1,2-Dichloroethene ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 0-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | Chloroethane | ND | ND | ug/l | NC | 20 |
| cis-1,2-Dichloroethene ND ND ug/l NC 20 Trichloroethene ND ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 o-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | 1,1-Dichloroethene | ND | ND | ug/l | NC | 20 |
| Trichloroethene ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ND NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 0-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | trans-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| 1,2-Dichlorobenzene ND ND ug/l NC 20 1,3-Dichlorobenzene ND ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 o-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | cis-1,2-Dichloroethene | ND | ND | ug/l | NC | 20 |
| 1,3-Dichlorobenzene ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 o-Xylene ND ND ND ug/l NC 20 1,1-Dichloropropene ND ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | Trichloroethene | ND | ND | ug/l | NC | 20 |
| 1,4-Dichlorobenzene ND ND ug/l NC 20 Styrene ND ND ND ug/l NC 20 o-Xylene ND ND ug/l NC 20 1,1-Dichloropropene ND ND ug/l NC 20 2,2-Dichloropropane ND ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ND ug/l NC 20 | 1,2-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| Styrene ND ND ug/l NC 20 o-Xylene ND ND ug/l NC 20 1,1-Dichloropropene ND ND ug/l NC 20 2,2-Dichloropropane ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20 | 1,3-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| o-Xylene ND ND ug/l NC 20 1,1-Dichloropropene ND ND ug/l NC 20 2,2-Dichloropropane ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20 | 1,4-Dichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,1-Dichloropropene ND ND ug/l NC 20 2,2-Dichloropropane ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20 | Styrene | ND | ND | ug/l | NC | 20 |
| 2,2-Dichloropropane ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20 | o-Xylene | ND | ND | ug/l | NC | 20 |
| 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20 | 1,1-Dichloropropene | ND | ND | ug/l | NC | 20 |
| | 2,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| 1.2.3-Trichloropropane ND ND ug/l NC 20 | 1,1,1,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| 7.55 - 7.51 - 1.52 - 2. | 1,2,3-Trichloropropane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L1004931

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|--------------------------------|-------------------|------------|-----------------|---------------------|
| /olatile Organics by GC/MS - Westborough Lab Sample | Associated sample(s): 01,05,09 | QC Batch ID: WG40 | 07461-6 QC | Sample: L100481 | 9-02 Client ID: DUP |
| Bromochloromethane | ND | ND | ug/l | NC | 20 |
| n-Butylbenzene | ND | ND | ug/l | NC | 20 |
| Dichlorodifluoromethane | ND | ND | ug/l | NC | 20 |
| Hexachlorobutadiene | ND | ND | ug/l | NC | 20 |
| Isopropylbenzene | ND | ND | ug/l | NC | 20 |
| p-Isopropyltoluene | ND | ND | ug/l | NC | 20 |
| Naphthalene | ND | ND | ug/l | NC | 20 |
| n-Propylbenzene | ND | ND | ug/l | NC | 20 |
| sec-Butylbenzene | ND | ND | ug/l | NC | 20 |
| tert-Butylbenzene | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| 1,3,5-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| Bromobenzene | ND | ND | ug/l | NC | 20 |
| o-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| p-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| Dibromomethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dibromoethane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L1004931

Report Date:

04/15/10

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|-----------------------------------|------------------|------------|-----------------------------|----------------------|
| Volatile Organics by GC/MS - Westborough L Sample | ab Associated sample(s): 01,05,09 | QC Batch ID: WG4 | 07461-6 QC | Sample: L10048 ² | 19-02 Client ID: DUP |
| 1,2-Dibromo-3-chloropropane | ND | ND | ug/l | NC | 20 |
| 1,3-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Methyl tert butyl ether | ND | ND | ug/l | NC | 20 |

| | | | | | Acceptance | |
|------------------------|-----------|-----------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 103 | | 104 | | 80-120 | |
| 4-Bromofluorobenzene | 92 | | 93 | | 80-120 | |

SEMIVOLATILES



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 Report Date: 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-01 Date Collected: 04/07/10 12:32

Client ID: GHC-5 Date Received: 04/07/10

Sample Location: WALPOLE, MA Field Prep: See Narrative
Matrix: Water Extraction Method: EPA 625

Analytical Method: 5,625 Extraction Date: 04/13/10 14:03

Analytical Date: 04/15/10 05:53
Analyst: PS

Parameter Result Qualifier Units RDL **Dilution Factor** Base/Neutral Extractables by GC/MS - Westborough Lab Acenaphthene ND ug/l 5.0 1 Benzidine ND ug/l 50 1 1,2,4-Trichlorobenzene ND 5.0 ug/l 1 Hexachlorobenzene ND 5.0 ug/l 1 ND Bis(2-chloroethyl)ether ug/l 5.0 1 ND 6.0 1 2-Chloronaphthalene ug/l 3,3'-Dichlorobenzidine ND ug/l 50 1 ND 6.0 2,4-Dinitrotoluene ug/l 1 2,6-Dinitrotoluene ND 5.0 ug/l 1 Azobenzene ND ug/l 5.0 1 Fluoranthene ND ug/l 5.0 1 4-Chlorophenyl phenyl ether ND 5.0 1 ug/l 4-Bromophenyl phenyl ether ND ug/l 5.0 1 ND Bis(2-chloroisopropyl)ether ug/l 5.0 1 Bis(2-chloroethoxy)methane ND ug/l 5.0 1 Hexachlorobutadiene ND ug/l 10 1 ND Hexachlorocyclopentadiene ug/l 30 1 Hexachloroethane ND 5.0 1 ug/l ND 5.0 Isophorone ug/l 1 Naphthalene ND 5.0 ug/l 1 Nitrobenzene ND ug/l 5.0 1 NDPA/DPA ND 15 1 ug/l n-Nitrosodi-n-propylamine ND ug/l 5.0 1 Bis(2-ethylhexyl)phthalate ND ug/l 5.0 1 Butyl benzyl phthalate ND ug/l 5.0 1 Di-n-butylphthalate ND ug/l 5.0 1 Di-n-octylphthalate ND ug/l 5.0 1 Diethyl phthalate ND ug/l 5.0 1 ND Dimethyl phthalate ug/l 5.0 1 ND 5.0 Benzo(a)anthracene ug/l 1



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-01 Date Collected: 04/07/10 12:32

Client ID: GHC-5 Date Received: 04/07/10
Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL **Dilution Factor Parameter** Result Qualifier Units Base/Neutral Extractables by GC/MS - Westborough Lab ND 5.0 Benzo(a)pyrene ug/l 1 ND 5.0 1 Benzo(b)fluoranthene ug/l Benzo(k)fluoranthene ND ug/l 5.0 1 5.0 Chrysene ND ug/l 1 Acenaphthylene ND 5.0 ug/l 1 Anthracene ND ug/l 5.0 1 Benzo(ghi)perylene ND ug/l 5.0 1 Fluorene ND ug/l 5.0 1 Phenanthrene ND ug/l 5.0 1 Dibenzo(a,h)anthracene ND ug/l 5.0 1 Indeno(1,2,3-cd)pyrene ND ug/l 7.0 1 Pyrene ND ug/l 5.0 1 Aniline ND ug/l 20 1 4-Chloroaniline ND ug/l 5.0 1 1-Methylnaphthalene ND ug/l 5.0 1 2-Nitroaniline ND 5.0 ug/l 1 3-Nitroaniline ND ug/l 5.0 1 ND 4-Nitroaniline ug/l 7.0 1 Dibenzofuran ND 5.0 1 ug/l 2-Methylnaphthalene ND ug/l 5.0 1 ND 50 n-Nitrosodimethylamine ug/l 1

| | | | Acceptance | |
|------------------|------------|-----------|------------|--|
| Surrogate | % Recovery | Qualifier | Criteria | |
| Nitrobenzene-d5 | 48 | | 23-120 | |
| 2-Fluorobiphenyl | 49 | | 15-120 | |
| 4-Terphenyl-d14 | 78 | | 33-120 | |



04/07/10

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-05 Date Collected: 04/07/10 13:39

Client ID: RIZ-2 Date Received:

Sample Location: WALPOLE, MA Field Prep: See Narrative Matrix: Water Extraction Method: EPA 625

Matrix: Water Extraction Method: EPA 625
Analytical Method: 5,625 Extraction Date: 04/13/10 14:03

Analytical Method: 5,625 Extraction Date: 04/13/10 14:03
Analytical Date: 04/15/10 06:18

Analyst: PS

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor |
|---|----------------|-----------|-------|-----|-----------------|
| Base/Neutral Extractables by GC/MS - We | estborough Lab | | | | |
| Acenaphthene | ND | | ug/l | 5.0 | 1 |
| Benzidine | ND | | ug/l | 50 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 5.0 | 1 |
| Hexachlorobenzene | ND | | ug/l | 5.0 | 1 |
| Bis(2-chloroethyl)ether | ND | | ug/l | 5.0 | 1 |
| 2-Chloronaphthalene | ND | | ug/l | 6.0 | 1 |
| 3,3'-Dichlorobenzidine | ND | | ug/l | 50 | 1 |
| 2,4-Dinitrotoluene | ND | | ug/l | 6.0 | 1 |
| 2,6-Dinitrotoluene | ND | | ug/l | 5.0 | 1 |
| Azobenzene | ND | | ug/l | 5.0 | 1 |
| Fluoranthene | ND | | ug/l | 5.0 | 1 |
| 4-Chlorophenyl phenyl ether | ND | | ug/l | 5.0 | 1 |
| 4-Bromophenyl phenyl ether | ND | | ug/l | 5.0 | 1 |
| Bis(2-chloroisopropyl)ether | ND | | ug/l | 5.0 | 1 |
| Bis(2-chloroethoxy)methane | ND | | ug/l | 5.0 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 10 | 1 |
| Hexachlorocyclopentadiene | ND | | ug/l | 30 | 1 |
| Hexachloroethane | ND | | ug/l | 5.0 | 1 |
| Isophorone | ND | | ug/l | 5.0 | 1 |
| Naphthalene | ND | | ug/l | 5.0 | 1 |
| Nitrobenzene | ND | | ug/l | 5.0 | 1 |
| NDPA/DPA | ND | | ug/l | 15 | 1 |
| n-Nitrosodi-n-propylamine | ND | | ug/l | 5.0 | 1 |
| Bis(2-ethylhexyl)phthalate | ND | | ug/l | 5.0 | 1 |
| Butyl benzyl phthalate | ND | | ug/l | 5.0 | 1 |
| Di-n-butylphthalate | ND | | ug/l | 5.0 | 1 |
| Di-n-octylphthalate | ND | | ug/l | 5.0 | 1 |
| Diethyl phthalate | ND | | ug/l | 5.0 | 1 |
| Dimethyl phthalate | ND | | ug/l | 5.0 | 1 |
| Benzo(a)anthracene | ND | | ug/l | 5.0 | 1 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-05 Date Collected: 04/07/10 13:39

Client ID: RIZ-2 Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL **Dilution Factor Parameter** Result Qualifier Units Base/Neutral Extractables by GC/MS - Westborough Lab ND 5.0 Benzo(a)pyrene ug/l 1 ND 5.0 1 Benzo(b)fluoranthene ug/l Benzo(k)fluoranthene ND ug/l 5.0 1 5.0 Chrysene ND ug/l 1 Acenaphthylene ND 5.0 ug/l 1 Anthracene ND ug/l 5.0 1 Benzo(ghi)perylene ND ug/l 5.0 1 Fluorene ND ug/l 5.0 1 Phenanthrene ND ug/l 5.0 1 Dibenzo(a,h)anthracene ND ug/l 5.0 1 Indeno(1,2,3-cd)pyrene ND ug/l 7.0 1 Pyrene ND ug/l 5.0 1 Aniline ND ug/l 20 1 4-Chloroaniline ND ug/l 5.0 1 1-Methylnaphthalene ND ug/l 5.0 1 2-Nitroaniline ND 5.0 ug/l 1 3-Nitroaniline ND ug/l 5.0 1 ND 4-Nitroaniline ug/l 7.0 1 Dibenzofuran ND 5.0 1 ug/l 2-Methylnaphthalene ND ug/l 5.0 1

| | | | Acceptance | |
|------------------|------------|-----------|------------|--|
| Surrogate | % Recovery | Qualifier | Criteria | |
| Nitrobenzene-d5 | 56 | | 23-120 | |
| 2-Fluorobiphenyl | 59 | | 15-120 | |
| 4-Terphenyl-d14 | 81 | | 33-120 | |

ND



50

1

ug/l

n-Nitrosodimethylamine

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09 Date Collected: 04/07/10 14:19

Client ID: RIZ-10 Date Received: 04/07/10

Sample Location: WALPOLE, MA Field Prep: See Narrative Matrix: Water Extraction Method: EPA 625

Analytical Method: 5,625 Extraction Date: 04/13/10 14:03

Analytical Date: 04/14/10 16:59

Analyst: PS

| Parameter | Result | Qualifier U | nits RDL | Dilution Factor |
|--------------------------------------|-----------------|-------------|----------|-----------------|
| Base/Neutral Extractables by GC/MS - | Westborough Lab | | | |
| Acenaphthene | ND | u | g/l 5.0 | 1 |
| Benzidine | ND | u | g/l 50 | 1 |
| 1,2,4-Trichlorobenzene | ND | u | g/l 5.0 | 1 |
| Hexachlorobenzene | ND | u | g/l 5.0 | 1 |
| Bis(2-chloroethyl)ether | ND | u | g/l 5.0 | 1 |
| 2-Chloronaphthalene | ND | u | g/l 6.0 | 1 |
| 3,3'-Dichlorobenzidine | ND | u | g/l 50 | 1 |
| 2,4-Dinitrotoluene | ND | u | g/l 6.0 | 1 |
| 2,6-Dinitrotoluene | ND | u | g/l 5.0 | 1 |
| Azobenzene | ND | u | g/l 5.0 | 1 |
| Fluoranthene | ND | u | g/l 5.0 | 1 |
| 4-Chlorophenyl phenyl ether | ND | u | g/l 5.0 | 1 |
| 4-Bromophenyl phenyl ether | ND | u | g/l 5.0 | 1 |
| Bis(2-chloroisopropyl)ether | ND | u | g/l 5.0 | 1 |
| Bis(2-chloroethoxy)methane | ND | u | g/l 5.0 | 1 |
| Hexachlorobutadiene | ND | u | g/l 10 | 1 |
| Hexachlorocyclopentadiene | ND | u | g/l 30 | 1 |
| Hexachloroethane | ND | u | g/l 5.0 | 1 |
| Isophorone | ND | u | g/l 5.0 | 1 |
| Naphthalene | ND | u | g/l 5.0 | 1 |
| Nitrobenzene | ND | u | g/l 5.0 | 1 |
| NDPA/DPA | ND | u | g/l 15 | 1 |
| n-Nitrosodi-n-propylamine | ND | u | g/l 5.0 | 1 |
| Bis(2-ethylhexyl)phthalate | ND | u | g/l 5.0 | 1 |
| Butyl benzyl phthalate | ND | u | g/l 5.0 | 1 |
| Di-n-butylphthalate | ND | u | g/l 5.0 | 1 |
| Di-n-octylphthalate | ND | u | g/l 5.0 | 1 |
| Diethyl phthalate | ND | u | g/l 5.0 | 1 |
| Dimethyl phthalate | ND | u | g/l 5.0 | 1 |
| Benzo(a)anthracene | ND | u | g/l 5.0 | 1 |
| | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09 Date Collected: 04/07/10 14:19

Client ID: RIZ-10 Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

RDL **Dilution Factor Parameter** Result Qualifier Units Base/Neutral Extractables by GC/MS - Westborough Lab ND 5.0 Benzo(a)pyrene ug/l 1 ND 5.0 1 Benzo(b)fluoranthene ug/l Benzo(k)fluoranthene ND ug/l 5.0 1 5.0 Chrysene ND ug/l 1 Acenaphthylene ND 5.0 ug/l 1 Anthracene ND ug/l 5.0 1 Benzo(ghi)perylene ND ug/l 5.0 1 Fluorene ND ug/l 5.0 1 Phenanthrene ND ug/l 5.0 1 Dibenzo(a,h)anthracene ND ug/l 5.0 1 Indeno(1,2,3-cd)pyrene ND ug/l 7.0 1 Pyrene ND ug/l 5.0 1 Aniline ND ug/l 20 1 4-Chloroaniline ND ug/l 5.0 1 1-Methylnaphthalene ND ug/l 5.0 1 2-Nitroaniline ND 5.0 ug/l 1 3-Nitroaniline ND ug/l 5.0 1 ND 4-Nitroaniline ug/l 7.0 1 Dibenzofuran ND 5.0 1 ug/l 2-Methylnaphthalene ND ug/l 5.0 1 ND 50 n-Nitrosodimethylamine ug/l 1

| | Acceptance | | | | | | | | |
|------------------|------------|-----------|----------|--|--|--|--|--|--|
| Surrogate | % Recovery | Qualifier | Criteria | | | | | | |
| Nitrobenzene-d5 | 52 | | 23-120 | | | | | | |
| 2-Fluorobiphenyl | 59 | | 15-120 | | | | | | |
| 4-Terphenyl-d14 | 77 | | 33-120 | | | | | | |

Extraction Method: EPA 625

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

Analytical Method: 5,625

Analytical Date: 04/14/10 15:44 Extraction Date: 04/13/10 14:03

Analyst: PS

| Parameter | Result | Qualifier | Units | RDL | | |
|-----------------------------------|-------------|-------------------|-----------|----------|--------|------------|
| Base/Neutral Extractables by GC/I | MS - Westbo | prough Lab for sa | ample(s): | 01,05,09 | Batch: | WG407902-1 |
| Acenaphthene | ND | | ug/l | 5.0 | | |
| Benzidine | ND | | ug/l | 50 | | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 5.0 | | |
| Hexachlorobenzene | ND | | ug/l | 5.0 | | |
| Bis(2-chloroethyl)ether | ND | | ug/l | 5.0 | | |
| 2-Chloronaphthalene | ND | | ug/l | 6.0 | | |
| 3,3'-Dichlorobenzidine | ND | | ug/l | 50 | | |
| 2,4-Dinitrotoluene | ND | | ug/l | 6.0 | | |
| 2,6-Dinitrotoluene | ND | | ug/l | 5.0 | | |
| Azobenzene | ND | | ug/l | 5.0 | | |
| Fluoranthene | ND | | ug/l | 5.0 | | |
| 4-Chlorophenyl phenyl ether | ND | | ug/l | 5.0 | | |
| 4-Bromophenyl phenyl ether | ND | | ug/l | 5.0 | | |
| Bis(2-chloroisopropyl)ether | ND | | ug/l | 5.0 | | |
| Bis(2-chloroethoxy)methane | ND | | ug/l | 5.0 | | |
| Hexachlorobutadiene | ND | | ug/l | 10 | | |
| Hexachlorocyclopentadiene | ND | | ug/l | 30 | | |
| Hexachloroethane | ND | | ug/l | 5.0 | | |
| Isophorone | ND | | ug/l | 5.0 | | |
| Naphthalene | ND | | ug/l | 5.0 | | |
| Nitrobenzene | ND | | ug/l | 5.0 | | |
| NDPA/DPA | ND | | ug/l | 15 | | |
| n-Nitrosodi-n-propylamine | ND | | ug/l | 5.0 | | |
| Bis(2-ethylhexyl)phthalate | ND | | ug/l | 5.0 | | |
| Butyl benzyl phthalate | ND | | ug/l | 5.0 | | |
| Di-n-butylphthalate | ND | | ug/l | 5.0 | | |
| Di-n-octylphthalate | ND | | ug/l | 5.0 | | |
| Diethyl phthalate | ND | | ug/l | 5.0 | | |
| Dimethyl phthalate | ND | | ug/l | 5.0 | | |
| Benzo(a)anthracene | ND | | ug/l | 5.0 | | |
| Benzo(a)pyrene | ND | | ug/l | 5.0 | | |



Extraction Method: EPA 625

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

Analytical Method: 5,625

Analytical Date: 04/14/10 15:44 Extraction Date: 04/13/10 14:03

Analyst: PS

| Parameter | Result | Qualifier | Units | RDL | | |
|---------------------------------|--------------|---------------|--------------|----------|--------|------------|
| Base/Neutral Extractables by GC | /MS - Westbo | orough Lab fo | r sample(s): | 01,05,09 | Batch: | WG407902-1 |
| Benzo(b)fluoranthene | ND | | ug/l | 5.0 | | |
| Benzo(k)fluoranthene | ND | | ug/l | 5.0 | | |
| Chrysene | ND | | ug/l | 5.0 | | |
| Acenaphthylene | ND | | ug/l | 5.0 | | |
| Anthracene | ND | | ug/l | 5.0 | | |
| Benzo(ghi)perylene | ND | | ug/l | 5.0 | | |
| Fluorene | ND | | ug/l | 5.0 | | |
| Phenanthrene | ND | | ug/l | 5.0 | | |
| Dibenzo(a,h)anthracene | ND | | ug/l | 5.0 | | |
| Indeno(1,2,3-cd)pyrene | ND | | ug/l | 7.0 | | |
| Pyrene | ND | | ug/l | 5.0 | | |
| Aniline | ND | | ug/l | 20 | | |
| 4-Chloroaniline | ND | | ug/l | 5.0 | | |
| 1-Methylnaphthalene | ND | | ug/l | 5.0 | | |
| 2-Nitroaniline | ND | | ug/l | 5.0 | | |
| 3-Nitroaniline | ND | | ug/l | 5.0 | | |
| 4-Nitroaniline | ND | | ug/l | 7.0 | | |
| Dibenzofuran | ND | | ug/l | 5.0 | | |
| 2-Methylnaphthalene | ND | | ug/l | 5.0 | | |
| n-Nitrosodimethylamine | ND | | ug/l | 50 | | |

| | | Acceptance | | | | | | | |
|------------------|-----------|------------|----------|--|--|--|--|--|--|
| Surrogate | %Recovery | Qualifier | Criteria | | | | | | |
| NE L | | | 00.400 | | | | | | |
| Nitrobenzene-d5 | 63 | | 23-120 | | | | | | |
| 2-Fluorobiphenyl | 65 | | 15-120 | | | | | | |
| 4-Terphenyl-d14 | 90 | | 33-120 | | | | | | |



Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

Report Date: 04/15/10

| arameter | LCS %Recovery | LCSD Qual %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|---------------------------|-------------|---------------------|-----|------|------------|
| ase/Neutral Extractables by GC/MS - Westb | orough Lab As | sociated sample(s): 01,08 | 5,09 Batch: | WG407902-2 | | | |
| Acenaphthene | 70 | - | | 46-118 | - | | 30 |
| 1,2,4-Trichlorobenzene | 54 | - | | 39-98 | - | | 30 |
| 2-Chloronaphthalene | 77 | - | | 40-140 | - | | 30 |
| 2,4-Dinitrotoluene | 81 | - | | 24-96 | - | | 30 |
| 2,6-Dinitrotoluene | 72 | - | | 40-140 | - | | 30 |
| Fluoranthene | 88 | - | | 40-140 | - | | 30 |
| 4-Chlorophenyl phenyl ether | 73 | - | | 40-140 | - | | 30 |
| n-Nitrosodi-n-propylamine | 60 | - | | 41-116 | - | | 30 |
| Butyl benzyl phthalate | 82 | - | | 40-140 | - | | 30 |
| Anthracene | 83 | - | | 40-140 | - | | 30 |
| Pyrene | 81 | - | | 26-127 | - | | 30 |
| P-Chloro-M-Cresol | 75 | - | | 23-97 | - | | 30 |
| 2-Chlorophenol | 61 | - | | 27-123 | - | | 30 |
| 2-Nitrophenol | 66 | - | | 30-130 | - | | 30 |
| 4-Nitrophenol | 57 | - | | 10-80 | - | | 30 |
| 2,4-Dinitrophenol | 52 | - | | 20-130 | - | | 30 |
| Pentachlorophenol | 68 | - | | 9-103 | - | | 30 |
| Phenol | 31 | - | | 12-110 | - | | 30 |



Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Lab Number: L1004931

Project Number: 12700058 Report Date:

04/15/10

| | LCS | | LCSD | | %Recovery | | | |
|------------------|-----------|------|-----------|------|-----------|-----|------|------------|
| <u>Parameter</u> | %Recovery | Qual | %Recovery | Qual | Limits | RPD | Qual | RPD Limits |
| | | | | | | | ' | |

Base/Neutral Extractables by GC/MS - Westborough Lab Associated sample(s): 01,05,09 Batch: WG407902-2

| | LCS | | LCSD | | Acceptance |
|----------------------|-----------|------|-----------|------|------------|
| Surrogate | %Recovery | Qual | %Recovery | Qual | Criteria |
| 2-Fluorophenol | 41 | | | | 21-120 |
| Phenol-d6 | 29 | | | | 10-120 |
| Nitrobenzene-d5 | 56 | | | | 23-120 |
| 2-Fluorobiphenyl | 64 | | | | 15-120 |
| 2,4,6-Tribromophenol | 78 | | | | 10-120 |
| 4-Terphenyl-d14 | 79 | | | | 33-120 |



Matrix Spike Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

Report Date: 04/15/10

| Parameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery Qua | Recovery I Limits | RPD | Qual | RPD <u>Limit</u> s |
|---------------------------------------|------------------|--------------|---------------|-----------------|---------|--------------|----------------------|----------------------|----------|----------|-----------------------|
| Base/Neutral Extractables b MS Sample | y GC/MS - Wes | stborough La | ab Associated | d sample(s): 0 | 1,05,09 | QC Batc | h ID: WG407902-3 | QC Samp | le: L100 |)4952-01 | Client ID: |
| Acenaphthene | ND | 97.6 | 73 | 75 | | - | - | 46-118 | - | | 30 |
| 1,2,4-Trichlorobenzene | ND | 97.6 | 62 | 64 | | - | - | 39-98 | - | | 30 |
| 2-Chloronaphthalene | ND | 97.6 | 83 | 85 | | - | - | 40-140 | - | | 30 |
| 2,4-Dinitrotoluene | ND | 97.6 | 83 | 85 | | - | - | 24-96 | - | | 30 |
| 2,6-Dinitrotoluene | ND | 97.6 | 78 | 80 | | - | - | 40-140 | - | | 30 |
| Fluoranthene | ND | 97.6 | 86 | 88 | | - | - | 40-140 | - | | 30 |
| 4-Chlorophenyl phenyl ether | ND | 97.6 | 75 | 77 | | - | - | 40-140 | - | | 30 |
| n-Nitrosodi-n-propylamine | ND | 97.6 | 65 | 67 | | - | - | 41-116 | - | | 30 |
| Butyl benzyl phthalate | ND | 97.6 | 85 | 87 | | - | - | 40-140 | - | | 30 |
| Anthracene | ND | 97.6 | 79 | 81 | | - | - | 40-140 | - | | 30 |
| Pyrene | ND | 97.6 | 82 | 84 | | - | - | 26-127 | - | | 30 |
| P-Chloro-M-Cresol | ND | 97.6 | 81 | 83 | | - | - | 23-97 | - | | 30 |
| 2-Chlorophenol | ND | 97.6 | 67 | 69 | | - | - | 27-123 | - | | 30 |
| 2-Nitrophenol | ND | 97.6 | 72 | 74 | | - | - | 30-130 | - | | 30 |
| 4-Nitrophenol | ND | 97.6 | 77 | 79 | | - | - | 10-80 | - | | 30 |
| 2,4-Dinitrophenol | ND | 97.6 | ND | 64 | | - | - | 20-130 | - | | 30 |
| Pentachlorophenol | ND | 97.6 | 78 | 80 | | - | - | 9-103 | - | | 30 |
| Phenol | ND | 97.6 | 54 | 55 | | - | - | 12-110 | - | | 30 |
| | | | | | | | | | | | |



Matrix Spike Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 Lab Number:

L1004931

Report Date:

04/15/10

| | Native | MS | MS | MS | | MSD | MSD | | Recovery | | | RPD |
|-----------|--------|-------|-------|-----------|------|-------|-----------|------|----------|-----|------|--------|
| Parameter | Sample | Added | Found | %Recovery | Qual | Found | %Recovery | Qual | Limits | RPD | Qual | Limits |
| | | | | | | | | | | | | |

Base/Neutral Extractables by GC/MS - Westborough Lab Associated sample(s): 01,05,09 QC Batch ID: WG407902-3 QC Sample: L1004952-01 Client ID:

MS Sample

| | MS | ; | M | SD | Acceptance | |
|----------------------|------------|-----------|------------|-----------|------------|--|
| Surrogate | % Recovery | Qualifier | % Recovery | Qualifier | Criteria | |
| 2,4,6-Tribromophenol | 81 | | | | 10-120 | |
| 2-Fluorobiphenyl | 72 | | | | 15-120 | |
| 2-Fluorophenol | 64 | | | | 21-120 | |
| 4-Terphenyl-d14 | 85 | | | | 33-120 | |
| Nitrobenzene-d5 | 65 | | | | 23-120 | |
| Phenol-d6 | 55 | | | | 10-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931 **Report Date:** 04/15/10

| arameter | Native Sample | Duplicate Sample | Units | RPD | Qual RPD Limits |
|--|--------------------------|-----------------------|------------|--------|------------------------------|
| ase/Neutral Extractables by GC/MS - Westborough L JP Sample | ab Associated sample(s): | 01,05,09 QC Batch ID: | WG407902-4 | QC Sam | nple: L1004952-01 Client ID: |
| Acenaphthene | ND | ND | ug/l | NC | 30 |
| Benzidine | ND | ND | ug/l | NC | 30 |
| 1,2,4-Trichlorobenzene | ND | ND | ug/l | NC | 30 |
| Hexachlorobenzene | ND | ND | ug/l | NC | 30 |
| Bis(2-chloroethyl)ether | ND | ND | ug/l | NC | 30 |
| 2-Chloronaphthalene | ND | ND | ug/l | NC | 30 |
| 3,3'-Dichlorobenzidine | ND | ND | ug/l | NC | 30 |
| 2,4-Dinitrotoluene | ND | ND | ug/l | NC | 30 |
| 2,6-Dinitrotoluene | ND | ND | ug/l | NC | 30 |
| Azobenzene | ND | ND | ug/l | NC | 30 |
| Fluoranthene | ND | ND | ug/l | NC | 30 |
| 4-Chlorophenyl phenyl ether | ND | ND | ug/l | NC | 30 |
| 4-Bromophenyl phenyl ether | ND | ND | ug/l | NC | 30 |
| Bis(2-chloroisopropyl)ether | ND | ND | ug/l | NC | 30 |
| Bis(2-chloroethoxy)methane | ND | ND | ug/l | NC | 30 |
| Hexachlorobutadiene | ND | ND | ug/l | NC | 30 |
| Hexachlorocyclopentadiene | ND | ND | ug/l | NC | 30 |
| Hexachloroethane | ND | ND | ug/l | NC | 30 |
| Isophorone | ND | ND | ug/l | NC | 30 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931 **Report Date:** 04/15/10

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------|-----------------------|------------|---------|-----------------------------|
| Base/Neutral Extractables by GC/MS - Westborough Lab DUP Sample | Associated sample(s): | 01,05,09 QC Batch ID: | WG407902-4 | QC Samp | ble: L1004952-01 Client ID: |
| Naphthalene | ND | ND | ug/l | NC | 30 |
| Nitrobenzene | ND | ND | ug/l | NC | 30 |
| NDPA/DPA | ND | ND | ug/l | NC | 30 |
| n-Nitrosodi-n-propylamine | ND | ND | ug/l | NC | 30 |
| Bis(2-ethylhexyl)phthalate | 37 | 46 | ug/l | 22 | 30 |
| Butyl benzyl phthalate | ND | ND | ug/l | NC | 30 |
| Di-n-butylphthalate | ND | ND | ug/l | NC | 30 |
| Di-n-octylphthalate | ND | ND | ug/l | NC | 30 |
| Diethyl phthalate | ND | ND | ug/l | NC | 30 |
| Dimethyl phthalate | ND | ND | ug/l | NC | 30 |
| Benzo(a)anthracene | ND | ND | ug/l | NC | 30 |
| Benzo(a)pyrene | ND | ND | ug/l | NC | 30 |
| Benzo(b)fluoranthene | ND | ND | ug/l | NC | 30 |
| Benzo(k)fluoranthene | ND | ND | ug/l | NC | 30 |
| Chrysene | ND | ND | ug/l | NC | 30 |
| Acenaphthylene | ND | ND | ug/l | NC | 30 |
| Anthracene | ND | ND | ug/l | NC | 30 |
| Benzo(ghi)perylene | ND | ND | ug/l | NC | 30 |
| Fluorene | ND | ND | ug/l | NC | 30 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

Report Date: 04/15/10

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|--------------------------|-----------------------|------------|---------|----------------------------|
| Base/Neutral Extractables by GC/MS - Westborough La DUP Sample | ab Associated sample(s): | 01,05,09 QC Batch ID: | WG407902-4 | QC Samp | le: L1004952-01 Client ID: |
| Phenanthrene | ND | ND | ug/l | NC | 30 |
| Dibenzo(a,h)anthracene | ND | ND | ug/l | NC | 30 |
| Indeno(1,2,3-cd)pyrene | ND | ND | ug/l | NC | 30 |
| Pyrene | ND | ND | ug/l | NC | 30 |
| Aniline | ND | ND | ug/l | NC | 30 |
| 4-Chloroaniline | ND | ND | ug/l | NC | 30 |
| 1-Methylnaphthalene | ND | ND | ug/l | NC | 30 |
| 2-Nitroaniline | ND | ND | ug/l | NC | 30 |
| 3-Nitroaniline | ND | ND | ug/l | NC | 30 |
| 4-Nitroaniline | ND | ND | ug/l | NC | 30 |
| Dibenzofuran | ND | ND | ug/l | NC | 30 |
| 2-Methylnaphthalene | ND | ND | ug/l | NC | 30 |
| n-Nitrosodimethylamine | ND | ND | ug/l | NC | 30 |

| | | | Acceptance | |
|------------------|-----------|---------------------|--------------------|--|
| Surrogate | %Recovery | Qualifier %Recovery | Qualifier Criteria | |
| Nitrobenzene-d5 | 58 | 71 | 23-120 | |
| 2-Fluorobiphenyl | 58 | 74 | 15-120 | |
| 4-Terphenyl-d14 | 81 | 93 | 33-120 | |



Lab Number: **Project Name:** WALPOLE PARK SOUTH L1004931

04/15/10 **Project Number:** Report Date: 12700058

Parameter Native Sample Duplicate Sample Units RPD **RPD Limits** Base/Neutral Extractables by GC/MS - Westborough Lab Associated sample(s): 01,05,09 QC Batch ID: WG407902-4 QC Sample: L1004952-01 Client ID:

DUP Sample

Acceptance Surrogate Criteria %Recovery Qualifier %Recovery Qualifier



METALS



64,7470A

60,6010B

60,6010B

ΕZ

ΑI

ΑI

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-01

Client ID: GHC-5

Sample Location: WALPOLE, MA

ND

ND

ND

Matrix: Water

Mercury, Dissolved

Selenium, Dissolved

Silver, Dissolved

Date Collected: 04/07/10 12:32

04/08/10 17:10 04/09/10 11:53 EPA 7470A

04/08/10 09:35 04/09/10 10:57 EPA 3005A

04/08/10 09:35 04/09/10 10:57 EPA 3005A

Date Received: 04/07/10

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|---------------------|----------|-------------|-------|-------|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Total Metals - | Westbord | ough Lab | | | | | | | | |
| Sodium, Total | 59 | | mg/l | 2.0 | 1 | 04/08/10 10:00 | 04/09/10 13:39 | EPA 3005A | 60,6010B | Al |
| MCP Dissolved Met | | tborough La | | | | | | | | |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | 1 | 04/08/10 09:35 | 04/09/10 10:57 | EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.171 | | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 10:57 | EPA 3005A | 60,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | 1 | 04/08/10 09:35 | 04/09/10 10:57 | EPA 3005A | 60,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | 1 | 04/08/10 09:35 | 04/09/10 10:57 | EPA 3005A | 60,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 10:57 | EPA 3005A | 60,6010B | Al |

1

1

1

0.0002

0.010

0.007

mg/l

mg/l

mg/l



04/07/10 13:39

04/07/10

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Date Collected:

Date Received:

Lab ID: L1004931-05

Client ID: RIZ-2

Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|--------------------|----------|-----------|-------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Total Metals - | Westbord | ough Lab | | | | | | | | |
| Sodium, Total | 68 | | mg/l | 2.0 | 1 | 04/08/10 10:00 | 04/09/10 13:46 | EPA 3005A | 60,6010B | Al |

| MCP Dissolved Me | etals - Westbor | ough Lab | | | | | | | |
|---------------------|-----------------|----------|--------|---|----------------|----------------|-----------|----------|----|
| Arsenic, Dissolved | ND | mg/l | 0.005 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Barium, Dissolved | 0.090 | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Cadmium, Dissolved | ND | mg/l | 0.004 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Chromium, Dissolved | ND | mg/l | 0.01 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Lead, Dissolved | ND | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 04/08/10 17:10 | 04/09/10 11:55 | EPA 7470A | 64,7470A | EZ |
| Selenium, Dissolved | ND | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| Silver, Dissolved | ND | mg/l | 0.007 | 1 | 04/08/10 09:35 | 04/09/10 11:03 | EPA 3005A | 60,6010B | AI |
| | | | | | | | | | |



60,6010B

60,6010B

60,6010B

60,6010B

60,6010B

64,7470A

60,6010B

60,6010B

ΑI

ΑI

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ΕZ

ΑI

ΑI

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09

Client ID: RIZ-10

Sample Location: WALPOLE, MA

ND

ND

ND

ND

ND

ND

ND

0.030

Matrix: Water

Arsenic, Dissolved

Barium, Dissolved

Cadmium, Dissolved

Chromium, Dissolved

Lead, Dissolved

Mercury, Dissolved

Selenium, Dissolved

Silver, Dissolved

Date Collected: 04/07/10 14:19

Date Received: 04/07/10

04/08/10 09:35 04/09/10 11:06 EPA 3005A

04/08/10 17:10 04/09/10 11:56 EPA 7470A

04/08/10 09:35 04/09/10 11:06 EPA 3005A

04/08/10 09:35 04/09/10 11:06 EPA 3005A

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|--------------------|------------|--------------|-------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Total Metals - | Westbord | ough Lab | | | | | | | | |
| Sodium, Total | 84 | | mg/l | 2.0 | 1 | 04/08/10 10:00 | 04/09/10 13:49 | EPA 3005A | 60,6010B | AI |
| | | | | | | | | | | |
| MCP Dissolved Met | tals - Wes | tborough Lab |) | | | | | | | |

1

1

1

1

1

1

1

1

0.005

0.010

0.004

0.01

0.010

0.0002

0.010

0.007

mg/l

mg/l

mg/l

mg/l

mg/l

mg/l

mg/l

mg/l



Project Name: WALPOLE PARK SOUTH **Lab Number:** L1004931

Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------|------------------------|----------|----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals | s - Westborough Lab fo | r sample | (s): 01, | 05,09 Bat | ch: WG40726 | 6-1 | | |
| Arsenic, Dissolved | ND | mg/l | 0.005 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | AI |
| Barium, Dissolved | ND | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | AI |
| Cadmium, Dissolved | ND | mg/l | 0.004 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | AI |
| Chromium, Dissolved | ND | mg/l | 0.01 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | AI |
| Lead, Dissolved | ND | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | AI |
| Selenium, Dissolved | ND | mg/l | 0.010 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | Al |
| Silver, Dissolved | ND | mg/l | 0.007 | 1 | 04/08/10 09:35 | 04/09/10 10:47 | 60,6010B | Al |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Q | ualifier Units | | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|------------------|------------------|------------------|----------|--------------------|------------------|------------------|----------------------|----|
| MCP Total Metals | - Westborough La | b for sample(s): | 01,05,09 | Batch: | WG407276-1 | | | |
| Sodium, Total | ND | mg/l | 2.0 | 1 | 04/08/10 10:00 | 04/09/10 13:19 | 60,6010B | Al |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|---------------------|-------------------------|----------|-----------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metal | ls - Westborough Lab fo | r sample | e(s): 01, | 05,09 Bat | ch: WG40734 | 5-1 | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | 1 | 04/08/10 17:10 | 04/09/10 11:48 | 64,7470A | EZ |

Prep Information

Digestion Method: EPA 7470A



Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1

L1004931

Report Date:

04/15/10

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|------------|-------------------|-------------|---------------------|-----|------|------------|
| MCP Dissolved Metals - Westborough Lab | Associated samp | ole(s): 01 | ,05,09 Batch | WG407266-2 | 2 WG407266-3 | | | |
| Arsenic, Dissolved | 112 | | 113 | | 80-120 | 1 | | 20 |
| Barium, Dissolved | 100 | | 100 | | 80-120 | 0 | | 20 |
| Cadmium, Dissolved | 113 | | 113 | | 80-120 | 0 | | 20 |
| Chromium, Dissolved | 95 | | 100 | | 80-120 | 5 | | 20 |
| Lead, Dissolved | 107 | | 107 | | 80-120 | 0 | | 20 |
| Selenium, Dissolved | 115 | | 115 | | 80-120 | 0 | | 20 |
| Silver, Dissolved | 97 | | 98 | | 80-120 | 1 | | 20 |
| MCP Total Metals - Westborough Lab Asso | ciated sample(s) | : 01,05,0 | 9 Batch: W | G407276-2 W | G407276-3 | | | |
| Sodium, Total | 98 | | 100 | | 80-120 | 2 | | 20 |
| MCP Dissolved Metals - Westborough Lab | Associated samp | ole(s): 01 | ,05,09 Batch | WG407345-2 | 2 WG407345-3 | | | |
| Mercury, Dissolved | 115 | | 115 | | 80-120 | 0 | | 20 |



INORGANICS & MISCELLANEOUS



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 Report Date: 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-01 Date Collected: 04/07/10 12:32

Client ID: GHC-5 Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|------------------------|---------------|-----------|-------|------|--------------------|------------------|------------------|----------------------|---------|
| General Chemistry - We | stborough Lab | | | | | | | | |
| Nitrogen, Nitrite | ND | | mg/l | 0.05 | 1 | - | 04/08/10 19:49 | 30,4500NO3-F | DD |
| Nitrogen, Nitrate | 6.5 | | mg/l | 0.10 | 1 | - | 04/08/10 19:49 | 30,4500NO3-F | DD |
| Oil & Grease, Hem-Grav | ND | | mg/l | 4.0 | 1 | 04/08/10 11:00 | 04/09/10 09:15 | 74,1664A | JO |
| Phenolics, Total | ND | | mg/l | 0.15 | 5 | - | 04/10/10 21:31 | 1,9065 | TH |



L1004931

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: Report Date: 04/15/10 12700058

SAMPLE RESULTS

Lab ID: Date Collected: L1004931-02 04/07/10 12:56

GHC-5-A Client ID: Date Received: 04/07/10 WALPOLE, MA Not Specified Sample Location: Field Prep:

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|--------------------------|-----------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis | s - Westborough | Lab | | | | | | | |
| Coliform Fecal (MF) | ND | | col/100ml | 10 | 10 | _ | 04/07/10 20:35 | 30 9222D | .IT |



L1004931

04/15/10

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 **Report Date:**

SAMPLE RESULTS

Lab ID: L1004931-03

GHC-5-B Client ID: WALPOLE, MA Sample Location:

Matrix: Water Date Collected: 04/07/10 12:57

Date Received: 04/07/10

Lab Number:

Not Specified Field Prep:

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------------|-------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis - | Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 10 | 10 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH

Lab Number:

L1004931

Project Number: 12700058

Report Date:

04/15/10

SAMPLE RESULTS

Lab ID:

L1004931-04

Client ID:

GHC-5-C

Sample Location: WALPOLE, MA

Matrix:

Water

Date Collected:

04/07/10 12:58

Date Received:

04/07/10

Field Prep:

Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|--------------------------|---------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis | - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 10 | 10 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 Report Date: 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-05

Client ID: RIZ-2

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 04/07/10 13:39

Date Received: 04/07/10

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|------------------------|---------------|-----------|-------|------|--------------------|------------------|------------------|----------------------|---------|
| General Chemistry - We | stborough Lab | | | | | | | | |
| Nitrogen, Nitrite | ND | | mg/l | 0.05 | 1 | - | 04/08/10 19:50 | 30,4500NO3-F | DD |
| Nitrogen, Nitrate | 0.47 | | mg/l | 0.10 | 1 | - | 04/08/10 19:50 | 30,4500NO3-F | DD |
| Oil & Grease, Hem-Grav | ND | | mg/l | 4.0 | 1 | 04/08/10 11:00 | 04/09/10 09:15 | 74,1664A | JO |
| Phenolics, Total | ND | | mg/l | 0.03 | 1 | - | 04/10/10 21:32 | 1,9065 | TH |



L1004931

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-06

RIZ-2-A Client ID: Sample Location: WALPOLE, MA

Matrix: Water Date Collected: 04/07/10 13:55

Date Received: 04/07/10 Field Prep:

Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|--------------------------|-----------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis | s - Westborough | ı Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 2.0 | 2 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH

Lab Number:

L1004931

Project Number: 12700058

Report Date:

04/15/10

SAMPLE RESULTS

Lab ID: L1004931-07

Client ID: RIZ-2-B Sample Location: WALPOLE, MA

Matrix: Water

Date Collected:

04/07/10 13:56

Date Received: Field Prep:

04/07/10

Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------------|---------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis - | - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 2.0 | 2 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH

Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-08

RIZ-2-C Client ID: Sample Location: WALPOLE, MA

Matrix: Water Date Collected: 04/07/10 13:57

Date Received: 04/07/10

Not Specified Field Prep:

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------------|---------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis - | - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 2.0 | 2 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-09 Date Collected: 04/07/10 14:19

Client ID: RIZ-10 Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|------------------------|---------------|-----------|-------|------|--------------------|------------------|------------------|----------------------|---------|
| General Chemistry - We | stborough Lab | | | | | | | | |
| Nitrogen, Nitrite | ND | | mg/l | 0.05 | 1 | - | 04/08/10 19:51 | 30,4500NO3-F | DD |
| Nitrogen, Nitrate | 1.0 | | mg/l | 0.10 | 1 | - | 04/08/10 19:51 | 30,4500NO3-F | DD |
| Oil & Grease, Hem-Grav | ND | | mg/l | 4.4 | 1.1 | 04/08/10 11:00 | 04/09/10 09:15 | 74,1664A | JO |
| Phenolics, Total | ND | | mg/l | 0.15 | 5 | - | 04/10/10 21:33 | 1,9065 | TH |



L1004931

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-10 Date Collected: 04/07/10 14:32

Client ID: RIZ-10-A Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: Not Specified

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|------------------------|------------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analys | is - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 10 | 10 | - | 04/07/10 20:35 | 30,9222D | JT |



L1004931

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058 **Report Date:** 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-11

RIZ-10-B Client ID: WALPOLE, MA Sample Location:

Matrix: Water Date Collected: 04/07/10 14:33

Date Received: 04/07/10 Field Prep:

Lab Number:

Not Specified

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|--------------------------|---------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis | - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 10 | 10 | - | 04/07/10 20:35 | 30,9222D | JT |



L1004931

04/07/10 14:34

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058 Report Date: 04/15/10

SAMPLE RESULTS

Lab ID: L1004931-12 Date Collected:

Client ID: RIZ-10-C Date Received: 04/07/10 Sample Location: WALPOLE, MA Field Prep: Not Specified

Matrix: Water

| Parameter | Result | Qualifier | Units | RDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|--------------------------|-----------------|-----------|-----------|-----|--------------------|------------------|------------------|----------------------|---------|
| Microbiological Analysis | s - Westborough | Lab | | | | | | | |
| Coliform, Fecal (MF) | ND | | col/100ml | 10 | 10 | - | 04/07/10 20:35 | 30,9222D | JT |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

Method Blank Analysis Batch Quality Control

| Parameter | Result Q | ualifier Unit | s RDI | Dilution Factor | | Date repared | Date Analyzed | Analytical Method | Analyst |
|------------------------|-------------------|---------------|-------------|-----------------|--------|-----------------|------------------|----------------------|---------|
| Microbiological Analy | sis - Westborough | Lab for samp | ole(s): 02- | 04,06-08 | ,10-12 | Batch: WO | G407183-1 | | |
| Coliform, Fecal (MF) | ND | col/1 | 00ml 1.0 |) 1 | | - | 04/07/10 20:35 | 30,9222D | JT |
| General Chemistry - ' | Westborough Lab | for sample(s) | : 01,05,09 | Batch: | WG407 | 7254-2 | | | |
| Oil & Grease, Hem-Grav | ND | m | g/l 4.0 |) 1 | 04/ | 08/10 11:00 | 04/09/10 09:15 | 74,1664A | JO |
| General Chemistry - ' | Westborough Lab | for sample(s) | : 01,05,09 | Batch: | WG407 | 7333-2 | | | |
| Nitrogen, Nitrate | ND | m | g/l 0.1 | 0 1 | | - | 04/08/10 19:40 | 30,4500NO3-F | DD |
| General Chemistry - ' | Westborough Lab | for sample(s) | : 01,05,09 | Batch: | WG407 | 7334-2 | | | |
| Nitrogen, Nitrite | ND | m | g/l 0.0 | 5 1 | | - | 04/08/10 19:42 | 30,4500NO3-F | = DD |
| General Chemistry - ' | Westborough Lab | for sample(s) | : 01,05,09 | Batch: | WG407 | 7584-1 | | | |
| Phenolics, Total | ND | m | g/l 0.0 | 3 1 | | - | 04/10/10 21:27 | 1,9065 | TH |



Lab Control Sample Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

ab Number: L1004931

Report Date: 04/15/10

| Parameter | LCS %Recovery | Qual | LCSD %Recover | 'Y Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|-------------------------------------|-----------------------|----------|------------------|------------|---------------------|-----|------|------------|
| General Chemistry - Westborough Lab | Associated sample(s): | 01,05,09 | Batch: | WG407254-1 | | | | |
| Oil & Grease, Hem-Grav | 108 | | - | | 78-114 | - | | 18 |
| General Chemistry - Westborough Lab | Associated sample(s): | 01,05,09 | Batch: | WG407333-1 | | | | |
| Nitrogen, Nitrate | 98 | | - | | 90-110 | - | | |
| General Chemistry - Westborough Lab | Associated sample(s): | 01,05,09 | Batch: | WG407334-1 | | | | |
| Nitrogen, Nitrite | 100 | | - | | 90-110 | - | | |
| General Chemistry - Westborough Lab | Associated sample(s): | 01,05,09 | Batch: | WG407584-2 | | | | |
| Phenolics, Total | 101 | | - | | 70-130 | - | | |



Matrix Spike Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931

Report Date: 04/15/10

| Parameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Qual Found | MSD %Recovery | Recover Qual Limits | y RPD | RPD Qual Limits |
|------------------------------|------------------|-------------|----------------|-----------------|-------------------|------------------|---------------------|------------|--------------------|
| General Chemistry - Westboro | ough Lab Assoc | ciated samp | ole(s): 01,05, | 09 QC Batc | h ID: WG407254-3 | QC Sample | e: L1004815-02 | Client ID: | MS Sample |
| Oil & Grease, Hem-Grav | ND | 43 | 38 | 89 | - | - | 78-114 | - | 18 |
| General Chemistry - Westboro | ough Lab Assoc | ciated samp | le(s): 01,05, | 09 QC Batc | h ID: WG407333-3 | QC Sample | e: L1004954-08 | Client ID: | MS Sample |
| Nitrogen, Nitrate | 0.81 | 4 | 4.7 | 97 | - | - | 83-113 | - | 17 |
| General Chemistry - Westboro | ough Lab Assoc | ciated samp | le(s): 01,05, | 09 QC Batc | h ID: WG407334-3 | QC Sample | e: L1004931-01 | Client ID: | GHC-5 |
| Nitrogen, Nitrite | ND | 4 | 4.0 | 100 | - | - | 80-120 | - | 20 |
| General Chemistry - Westboro | ough Lab Assoc | ciated samp | le(s): 01,05, | 09 QC Batc | h ID: WG407584-3 | QC Sample | e: L1004931-05 | Client ID: | RIZ-2 |
| Phenolics, Total | ND | 0.8 | 0.73 | 91 | - | - | 70-130 | - | 20 |

Lab Duplicate Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number:

L1004931

Report Date:

04/15/10

| Parameter | Native Sample | Duplicate Sample | Units | RPD | Qual RPD Limits |
|---|---------------------|----------------------|------------|-------------|-----------------------|
| General Chemistry - Westborough Lab Associated samp | ole(s): 01,05,09 QC | Batch ID: WG407254-4 | QC Sample: | L1004835-03 | Client ID: DUP Sample |
| Oil & Grease, Hem-Grav | ND | ND | mg/l | NC | 18 |
| General Chemistry - Westborough Lab Associated samp | ole(s): 01,05,09 QC | Batch ID: WG407333-4 | QC Sample: | L1004954-08 | Client ID: DUP Sample |
| Nitrogen, Nitrate | 0.81 | 0.81 | mg/l | 0 | 17 |
| General Chemistry - Westborough Lab Associated samp | ole(s): 01,05,09 QC | Batch ID: WG407334-4 | QC Sample: | L1004931-01 | Client ID: GHC-5 |
| Nitrogen, Nitrite | ND | ND | mg/l | NC | 20 |
| General Chemistry - Westborough Lab Associated samp | ole(s): 01,05,09 QC | Batch ID: WG407584-4 | QC Sample: | L1004931-05 | Client ID: RIZ-2 |
| Phenolics, Total | ND | ND | mg/l | NC | 20 |



04151012:21

Project Name: WALPOLE PARK SOUTH Lab Number: L1004931

Project Number: 12700058 **Report Date:** 04/15/10

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal

A Absent VOA/VPH H2O Preservative Date: NA

B Absent

| Container Info | rmation | | | Temp | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|---|
| Container ID | Container Type | Cooler | рН | deg C | Pres | Seal | Analysis |
| L1004931-01A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-01B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-01C | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-01D | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-01E | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-01F | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-01G | Amber 1000ml H2SO4 preserved | Α | <2 | 5.9 | Υ | Absent | TPHENOL-9065(28) |
| L1004931-01H | Plastic 250ml HNO3 preserved | Α | <2 | 5.9 | Υ | Absent | MCP-NA-6010T(180) |
| L1004931-01I | Plastic 500ml unpreserved | Α | 7 | 5.9 | Υ | Absent | NO3-4500(2),NO2-4500NO3(2) |
| L1004931-01J | Plastic 500ml HNO3 preserved | A | <2 | 5.9 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SE-6010S(180),MCP-CD-6010S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180) |
| L1004931-02A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-02B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-03A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-03B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-04A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-04B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-05A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-05B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-05C | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-05D | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-05E | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-05F | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-05G | Amber 1000ml H2SO4 preserved | Α | <2 | 5.9 | Υ | Absent | TPHENOL-9065(28) |
| L1004931-05H | Plastic 250ml HNO3 preserved | Α | <2 | 5.9 | Υ | Absent | MCP-NA-6010T(180) |
| L1004931-05I | Plastic 500ml unpreserved | Α | 7 | 5.9 | Υ | Absent | NO3-4500(2),NO2-4500NO3(2) |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058

Lab Number: L1004931 **Report Date:** 04/15/10

| Container Info | ormation | | | Temp | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|---|
| Container ID | Container Type | Cooler | рΗ | deg C | Pres | Seal | Analysis |
| L1004931-05J | Plastic 500ml HNO3 preserved | Α | <2 | 5.9 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SE-6010S(180),MCP-CD-6010S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180) |
| L1004931-06A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-06B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-07A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-07B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-08A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-08B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-09A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-09B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 5.9 | Υ | Absent | 524.2(14) |
| L1004931-09C | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-09D | Amber 1000ml HCl preserved | В | N/A | 4 | Υ | Absent | OG-1664(28) |
| L1004931-09E | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-09F | Amber 1000ml unpreserved | В | 7 | 4 | Υ | Absent | BNEXT-625(7) |
| L1004931-09G | Amber 1000ml H2SO4 preserved | Α | <2 | 5.9 | Υ | Absent | TPHENOL-9065(28) |
| L1004931-09H | Plastic 250ml HNO3 preserved | Α | <2 | 5.9 | Υ | Absent | MCP-NA-6010T(180) |
| L1004931-09I | Plastic 500ml unpreserved | Α | 7 | 5.9 | Υ | Absent | NO3-4500(2),NO2-4500NO3(2) |
| L1004931-09J | Plastic 500ml HNO3 preserved | A | <2 | 5.9 | Y | Absent | MCP-AG-6010S(180),MCP-BA-6010S(180),MCP-SE-6010S(180),MCP-CD-6010S(180),MCP-7470S(28),MCP-CR-6010S(180),MCP-PB-6010S(180),MCP-AS-6010S(180) |
| L1004931-10A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-10B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-11A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-11B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-12A | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |
| L1004931-12B | Bacteria Cup Na2S2O3 preserved | Α | N/A | 5.9 | Υ | Absent | F-COLI-MF(.33) |



Project Name: WALPOLE PARK SOUTH Lab Number: L1004931
Project Number: 12700058 Report Date: 04/15/10

GLOSSARY

Acronyms

EPA · Environmental Protection Agency.

LCS Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD · Laboratory Control Sample Duplicate: Refer to LCS.

MS • Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD · Matrix Spike Sample Duplicate: Refer to MS.

NA · Not Applicable.

NC · Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

NI · Not Ignitable.

RDL - Reported Detection Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A Spectra identified as "Aldol Condensation Product".
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- ${f P}$ The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RDL. (Metals only.)
- ${f R}$ Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- J Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- **ND** Not detected at the reported detection limit (RDL) for the sample.

Report Format: Data Usability Report



Project Name:WALPOLE PARK SOUTHLab Number:L1004931Project Number:12700058Report Date:04/15/10

REFERENCES

- Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I IIIA, 1997.
- Methods for the Organic Chemical Analysis of Municipal and Industrial Wastewater. Appendix A, Part 136, 40 CFR (Code of Federal Regulations).
- Methods for the Determination of Organic Compounds in Drinking Water Supplement II. EPA/600/R-92/129, August 1992.
- 30 Standard Methods for the Examination of Water and Wastewater. APHA-AWWA-WPCF. 18th Edition. 1992.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). May 2004.
- Quality Assurance and Quality Control Requirements and Performance Standards for SW-846 Methods. MADEP BWSC. WSC-CAM-IIA (Revision 4), WSC-CAM-V C (Revision 2), WSC-CAM-IIIA (Revision 5). August 2004.
- Method 1664,Revision A: N-Hexane Extractable Material (HEM; Oil & Grease) and Silica Gel Treated N-Hexane Extractable Material (SGT-HEM; Non-polar Material) by Extraction and Gravimetry, EPA-821-R-98-002, February 1999.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Woods Hole Labs shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Woods Hole Labs.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised March 16, 2010 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. NELAP Accredited Solid Waste/Soil.

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate. Organic Parameters: Haloacetic Acids, Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB).)

Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Calcium Hardness, Silica, Sulfate, Sulfide, Ammonia, Kjeldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH.) Solid Waste/Soil (Inorganic Parameters: Lead in Paint, pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), Reactivity. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH, Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3.3'-Dichlorobenzidine, Phthalates, Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9221E, 9222B, 9222D, 9223B, EPA 180.1, 300.0, 353.2, SM2130B, 2320B, 4500Cl-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B,4500NO3-F, EPA 200.7, EPA 200.8, 245.1. Organic Parameters: 504.1, 524.2, SM 6251B.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, Lachat 10-107-06-1-B, SM2320B, 2340B, 2510B, 2540C, 2540D, 426C, 4500Cl-D, 4500Cl-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500P-B.5, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Drinking Water

Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl)

(EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Fluoride, Sulfate)

353.2 for: Nitrate-N, Nitrite-N; SM4500NO3-F, 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, 2320B, SM2540C. SM4500H-B.

Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics)

(504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), 314.0, 332.

Microbiology Parameters: SM9215B; ENZ. SUB. SM9223; MF-SM9222D

Non-Potable Water

Inorganic Parameters:, (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn)

(EPA 200.7 for: Al,Sb,As,Be,Cd,Cr,Co,Cu,Fe,Pb,Mn,Mo,Ni,Se,Aq,Sr,Ti,Tl, V,Zn,Ca,Mq,Na,K)

245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2540B, 2340B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Ammonia-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-B,C-Titr, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B,E, 5220D, EPA 410.4, SM 5210B, 5310C, 4500CL-D, EPA 1664, SM14 510AC, EPA 420, SM4500-CN-CE, SM2540D.

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics)

(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCBs-Water), EPA 625 for SVOC Acid Extractables and SVOC Base/Neutral Extractables, 600/4-81-045-PCB-Oil

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM6215B, 9222B, 9223B Colilert, EPA 200.7, 200.8, 245.2, 120.1, 300.0, 314.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 331.0. Organic Parameters: 504.1, 524.2, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 300.0, 350.1, 351.1, 353.2, 420.1, 1664A, SW-846 9010, 9030, 9040B, SM426C, SM2310B, 2540B, 2540D, 4500H+B, 4500NH3-H, 4500NH3-E, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 2320B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-117-07-1-B, LACHAT 10-107-06-1-B, LACHAT 10-107-04-1-J, LACHAT 10-117-07-1-A, SM4500CL-E, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3005A, 3015A, 3510C, 5030B, 8021B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 7.3.3.2, 7.3.4.2, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040, 9045C, 9050C, 1311, 3005A, 3050B, 3051A. Organic Parameters: SW-846 3540C, 3545, 3580A, 5030B, 5035, 8021B, 8260B, 8270C, 8330, 8151A, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 2540C, 2320B, 314.0, SM2120B, 2510B, 5310C, SM4500H-B, EPA 200.8, 245.2. Organic Parameters: 504.1, SM6251B, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.4, SM5220D, 4500Cl-D, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, SM9221CE, 9222D, 9221B, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.1, SM5210B, SW-846 3015, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 3510C, EPA 608, 624, 625, SW-846 5030B, 8021B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 9040B, 3005A, 6010B, 7196A, 5030B, 9010B, 9030B, 1030, 1311, 3050B, 3051, 7471A, 9014, 9012A, 9045C, 9050A, 9065. Organic Parameters: SW-846 8021B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 1311, 1312, 3540C, 3545, 3550B, 3580A, 5035L, 5035H, NJ OQA-QAM-025 Rev.7.)

New York Department of Health Certificate/Lab ID: 11148. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 9215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 314.0, 332.0, SM2320B, EPA 300.0, SM2120B, 4500CN-E, 4500F-C, 4500H-B, 4500NO3-F, 2540C, EPA 120.1, SM 2510B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, 5210B, EPA 410.4, SM5220D, 2310B-4a, 2320B, EPA 200.7, 300.0, LACHAT 10-117-07-1A or B, SM4500Cl-E, 4500F-C, SM15 426C, EPA 350.1, LACHAT 10-107-06-1-B, SM4500NH3-H, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-041-C, SM4500-NO3-F, 4500-NO2-B, 4500P-E, 2540C, 2540B, 2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, S\M3500Cr-D, EPA 245.1, 245.2, 7470A, SM2120B, SM4500-CN-E LACHAT 10-204-00-1-A, EPA 9040B, SM4500-HB, EPA 1664A, SM5310C, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, SM4500S-D, SM5540C, EPA 3005A, 3015. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, EPA 3510C, 5030B, 9010B, 9030B.)

Solid & Hazardous Waste (Inorganic Parameters: 1010, 1030, SW-846 Ch 7 Sec 7.3, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 1312, 3005A, 3050B, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8151A, 8330, 8082, 3540C, 3545, 3546, 3580, 5030B, 5035.)

North Carolina Department of the Environment and Natural Resources <u>Certificate/Lab ID</u>: 666. <u>Organic Parameters</u>: MA-EPH, MA-VPH.

Pennsylvania Department of Environmental Protection Certificate/Lab ID: 68-03671. *NELAP Accredited. Non-Potable Water* (Organic Parameters: EPA 3510C, 5030B, 625, 624. 608, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1010, 1030, 1311, 3050B, 3051, 6010B, EPA 7.3.3.2, EPA 7.3.4.2, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065. Organic Parameters: 3540C, 3545, 3580A, 5035, 8021B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Rhode Island Department of Health Certificate/Lab ID: LAO00065. *NELAP Accredited via NY-DOH.*Refer to MA-DEP Certificate for Potable and Non-Potable Water.
Refer to NY-DOH Certificate for Potable and Non-Potable Water.

Texas Commisson on Environmental Quality Certificate/Lab ID: T104704476-09-1. **NELAP Accredited.** Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664, 200.7, 200.8, 245.1, 245.2, 300.0, 350.1, 351.1, 353.2, 376.2, 410.4, 420.1, 6010, 6020, 7196, 7470, 9040, SM 2120B, 2310B, 2320B, 2510B, 2540B, 2540B, 2540D, 426C, 4500CL-E, 4500CN-E, 4500F-C, 4500H+B, 4500NH3-H, 4500NO2B, 4500P-E, 4500 S2⁻ D, 510C, 5210B, 5220D, 5310C, 5540C. Organic Parameters: EPA 608, 624, 625, 8081, 8082, 8151, 8260, 8270, 8330.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 9012, 9014, 9040, 9045, 9050, 9065.)

Utah Department of Health <u>Certificate/Lab ID</u>: AAMA. **NELAP Accredited.** *Non-Potable Water* (Inorganic Parameters: Chloride EPA 300.0)

Department of Defense Certificate/Lab ID: L2217.

Drinking Water (Inorganic Parameters: SM 4500H-B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: EPA 200.7, 200.8, 6010B, 6020, 245.1, 245.2, 7470A, 9040B, 300.0, 9251, 9038, 350.1, 353.2, 351.1, 314, 120.1, 9050A, 410.4, 9060, 1664, 420.1, LACHAT 10-107-06-1-B, SM 4500CN-E, 4500H-B, 4500CL-E, 4500F-BC, 4500SO4-E, 426C, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500NO2-B, 4500Norg-C, 4500PE, 2510B, 5540C, 5220D, 5310C, 2540B, 2540C, 2540D, 510C, 4500S2-AD, 3005A, 3015, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8330, 625, 8082, 8151A, 8081A, 3510C, 5030B.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 200.7, 6010B, 7471A, 9040B, 9045C, 9065, 420.1, 9012A, 6860, 1311, 1312, 3050B, 9030B, 3051, 9010B, 3540C, SM 510ABC, 4500CN-CE, 2540G, SW-846 7.3, Organic Parameters: EPA 8260B, 8270C, 8330, 8082, 8081A, 8151A, 3545, 3546, 3580, 5035.)

Analytes Not Accredited by NELAP

Certification is not available by NELAP for the following analytes: **EPA 8260B:** Freon-113, 1,2,4,5-Tetramethylbenzene, 4-Ethyltoluene. **EPA 8330A:** PETN, Picric Acid, Nitroglycerine, 2,6-DANT, 2,4-DANT. **EPA 8270C:** Methyl naphthalene, Dimethyl naphthalene, Total Methylnapthalenes, Total Dimethylnaphthalenes, 1,4-Diphenylhydrazine (Azobenzene). **EPA 625:** 4-Chloroaniline. **EPA 350.1** for Ammonia in a Soil matrix.

| FORM NO: 01-01 (rev. 18-Jan-2010) | T RCP? | IS YOUR BROIECT | PLEASE ANSWER QUESTIONS ABOVE | -04 bHC-5-C | -03 GHC-S-B | -07 bHC-5-A | . 4 | | | | | | #3101 PR-X | (Lab Use Only) Sample ID | ALPHA Lab ID | If MS is required, indicate in Sample Specific Comments which samples and what tests MS to be performed. (Note: All <i>CAM</i> methods for inorganic analyses require MS every 20 soil samples) | Alpha | Charlesh. 10m | | | inapan MA | Some Start S | Client: The Project Pr | Client Information Project Pro | TEL: 508-822-9300 | MANGETTI DI MA |
|-----------------------------------|--------------------------|--|-------------------------------------|-------------|-------------|-------------|--------|----------|------|----------|-----|--------|--|--------------------------|------------------------|---|--------------------------|---------------|--|--|---|--|--|--|-------------------------------|--|
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| | 1520 (held we + | Preservative WI - B N A | Container Type 46 A6 IN P | < | | | | | | 水 | × | × | X | Initials / (| Sampler's | ANA Verbly (6 Phenel | 127516 (264) (257) | | ☐ Yes ☐ No/ | DYes □ No s | | Mr. Thaces State /Fed Program | Regulatory Require | X ADEx | □FAX | Report Informatio |
| | By: Date/Time | ACI -N | 1 P P P | × | X | × | * | X | × | | | | | 1. ~ 4 | ~ 2 | SZX. Z NOZ NOZ NOZ NOZ NOZ NOZ NOZ NOZ NOZ NO | 1 | | Are CT RCP (Reasonable Confidence Protocols) Required? | Are MCP Analytical Methods Required? Is Matrix Spike (MS) Required on this SDG? (If yes see note in Comments) | MA MCP PRESUMPTIVE CERTAINTY CT REASONABLE CONFIDENCE PROTO | MCP Criteria RC | Requirements/Report Limits | eliverables (/c | | formation - Data Deliverables Billin |
| Cooligydiad ams. | 98967 (C-17C) (1-15) | pietely. Samples can not be logged in and turnaround time clock will not 8 | Please print clearly, legibly and c | | | | | | | | | | 77.4 | Sample Specific Comments | (Please specify below) | ■ Not needed □ Not needed □ Lab to do □ Preservation □ Lab to do | Filtration day | | Protocols) Required? | ? SDG?(If yes see note in Comment | ONABLE CONFIDENCE PROT | R(6W=1 | | AP | XSame as Client info PO #: | Billing Information |

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| FORM NO: 01-01 (rev. 18-Jan-2010) | MA MCP or CT RCP? | IS YOUR PROJECT | PLEASE ANSWER QUESTIONS ABOVE! | <u> </u> | | - 06 | | | | | | | E-21/1 50=1864 | (Lab Use Only) | AL DHA SAID | Other Project Specific Requirements/Comments/Detection Limits: If MS is required , indicate in Sample Specific Comments which samples and what tests MS to be performed. (Note: All CAM methods for inorganic analyses require MS every 20 soil samples) | ☐ These samples have been previously analyzed by Alpha | Email: A. J. A. | Fax: | Phone | 1.2 | Address: | Client: ************************************ | ation | పర | WESTBORO, MA MANSFIELD, MA | e C |
| | | | ABOVE! | 5 | 20 | 14 | | | | | | | | Sample ID | | ements/Comment beclific Comments which analyses require MS eve | analyzed by Alpha | 3 K | | | A | Pr | 14 P | | | | |
| | Refinquished By: | | | V 13 | 2 | | | | | | 3 | 3 (3 | 1/7/10 / | Date | Collection | s/Detection Limits: h samples and what tests ery 20 soil samples) | 1 1 1 1 me. 4/14/10 | | 1.5 | Turn-Around Time | ALPHA Quote #: | Project Manager: | Project #: | Project Location: | Project Name: | Project Information | CHAIN OF CUSTODY |
| | Date/Time | Preservative | Container Type | 1357 V 1 | 1356 | 1355 | 753 | <u>\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ </u> | 1350 | 1348 | 346 | 1344 | 1339 bw /4 | ime Matrix | Sample | MS to be performed. | | ☐ RUSH (only confirmed if pre-approved!) Time: | | | | | N. T. | 0 | | | PAGE OF _ |
| | 17 | /ative Mcl - 1 | Type H NAR | | | | | | | | × | × | * | Initials / O/ | 1 | ANAI Seas (66 Vartals (1) | Ys _i | S | □ Yes □ | DAYes □ No □ Yes □ No | 7 | State /Fed Pr | . ≺ | | □ FAX | Report Info | Date Rec'd in |
| | Received By: | N HCl - N | d | | | | × | <u>×</u> | X | * | | 2 2 | | | te 1 (0) | Phenol | 652 | 7 | | . 44 | ESUMPTIVE CERT | Program | Requirements/Report Limits | ☐ Add'l Deliverables | DEMAIL. | formation - Data Deliverables | in Lab: 4/7/10 |
| | Date/Time #17/10 /526 | 1 | P | <u> </u> | ∠ | ≺ | | | | | | | | 4 | 2/2000 2000 | RICES & COLO | net | / K | asonable Confidence | Are MCP Analytical Methods Required? Is Matrix Spike (MS) Required on this SI | AINTY CIT REAS | Criteria [2 | | ables | San | | |
| See reverse side. | | pletely, Samples in and turnaround | Please print clear | | | | | | | | | : | | Sample Specific Comments | (Please specify below) | ☐ Not needed ☐ Lab to do Preservation ☐ Lab to do | SAMPLE | | Are CT RCP (Reasonable Confidence Protocols) Required? | Are MCP Analytical Methods Required? Is Matrix Spike (MS) Required on this SDG?(If yes see note in Comments) | RESUMPTIVE CERTAINTY CT REASONABLE CONFIDENCE PROTO | 12 6 5 1 | A.A.A.A.A.A.A.A.A.A.A.A.A.A.A.A.A.A.A. | | Same as Client info PO#: | Billing Information | ALPHA Job#: U\C |
| Page | start until any ambiguittes are resolved. All samples submitted are subject to Alpha's Terms and Conditions. | pletely. Samples can not be logged in and turnaround time clock will not | Please print clearly, legibly and com- | ಬ | Ŋ | ىع | _ | | S | - | | \$J | ಲ | | y below) | | SAMPLE HANDLING | 0 1 | <u> </u> | e in Comments) | ENCE PROIO | | ender der State der State der State der State der State der State der State der State der State der State der | | # 25.00.00.00.00.00.00 | | 1004931 |
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| Ш | Sample Specific Comments | | de | \prec | (C) | 0 | | | Date Time | | Sample ID | Sam | (Lab Use Only) | (Lab l |
| | (Please specify below) | | 1/2 | 3 | | Ĵ/ | Sampler's | Sample | Collection | | | | ALPHA Lab ID | ALPH |
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| ئر ^ا <u>چ</u> | Filtration A. 2. | Low | 4.5 | | 16 | VAL | | | on Limits: | nts/Detection | nts/Comme | fic Requiremen | Other Project Specific Requirements/Comments/Detection Limits | Other |
| DLING T | SAMPLE HANDLING | 1 | _ | | ر که وی ک | rsis | | Time: | 0111111 | Date Due: 4/14/10 | zed by Alpha | en previously analyz | These samples have been previously analyzed by Alpha | ☐ The |
| | | te)s | | | | | -approved() | ly confirmed if pre- | ☐ RUSH (only confirmed if pre-approved!) | □ Standard | | | | Email: |
| omments) | Is Matrix Spike (MS) Required on this SDG? (If yes see note in Comments) Are CT RCP (Reasonable Confidence Protocols) Required? | Required on this Sable, Confidence P | Spike (MS) F | Is Matrix Are CT R | i | □ Yes | | | and Time | Turn-Around Time | | 7 | ¥ | Phone: |
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| 43) | ацрна Job #: ЦЮЦ93 | ALPHA | 011 | in Lab: 4/7/10 | | Date Rec'd | \f\{\partial \nabla \na | PAGE √3 | • | CUS | Z Q | CHAIN OF CUSTODY | | |
| | | 20 SS SE MICHAEL ST. 20 | | | | 100000000000000000000000000000000000000 | | | | | | | | |



ANALYTICAL REPORT

Lab Number: L1008812

Client: Tetra Tech Rizzo

1 Grant Street

Framingham, MA 01701-9005

ATTN: Ian Cannan Phone: (508) 903-2039

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Report Date: 06/21/10

Certifications & Approvals: MA (M-MA086), NY NELAC (11148), CT (PH-0574), NH (2003), NJ (MA935), RI (LAO00065), ME (MA0086), PA (Registration #68-03671), USDA (Permit #S-72578), US Army Corps of Engineers, Naval FESC.

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

| Alpha Sample ID | Client ID | Sample Location | Collection Date/Time |
|--------------------|---------------------|--------------------|-------------------------|
| L1008812-01 | MW-3 | WALPOLE, MA | 06/11/10 08:02 |
| L1008812-02 | RIZ-3 | WALPOLE, MA | 06/11/10 08:56 |
| L1008812-03 | MW-2 | WALPOLE, MA | 06/11/10 09:33 |
| L1008812-04 | GHC-6 | WALPOLE, MA | 06/11/10 10:08 |
| L1008812-05 | RIZ-9 | WALPOLE, MA | 06/11/10 11:07 |
| L1008812-06 | RIZ-10 | WALPOLE, MA | 06/11/10 12:30 |
| L1008812-07 | RIZ-8 | WALPOLE, MA | 06/11/10 13:40 |
| L1008812-08 | MW-9 | WALPOLE, MA | 06/11/10 14:10 |
| L1008812-09 | 20100611-TRIP BLANK | WALPOLE, MA | 06/11/10 00:00 |

Project Name: Lab Number: WALPOLE PARK SOUTH L1008812 **Project Number:** 12700058-003 **Report Date:**

06/21/10

MADEP MCP Response Action Analytical Report Certification

This form provides certifications for all samples performed by MCP methods. Please refer to the Sample Results and Container Information sections of this report for specification of MCP methods used for each analysis. The following questions pertain only to MCP **Analytical Methods.**

| A | Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature) in the field or laboratory, and prepared/analyzed within method holding times? | YES |
|----|---|-----|
| В | Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed? | NO |
| С | Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances? | YES |
|) | Does the laboratory report comply with all the reporting requirements specified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | YES |
| Ξa | VPH, EPH, and APH Methods only: Was each method conducted without significant modification(s)? (Refer to the individual method(s) for a list of significant modifications). | N/A |
| Ξb | APH and TO-15 Methods only: Was the complete analyte list reported for each method? | N/A |
| = | Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)? | YES |

| A res | sponse to questions G, H and I is required for "Presumptive Certainty" status | |
|-------|---|----|
| G | Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)? | NO |
| Н | Were all QC performance standards specified in the CAM protocol(s) achieved? | NO |
| I | Were results reported for the complete analyte list specified in the selected CAM protocol(s)? | NO |

For any questions answered "No", please refer to the case narrative section on the following page(s).

Please note that sample matrix information is located in the Sample Results section of this report.



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet all of the requirements of NELAC, for all NELAC accredited parameters. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

| For additional information, please contact Client Services at 800-624-9220. | |
|---|--|
| | |

MCP Related Narratives

Sample Receipt

The samples were Field Filtered for Dissolved Metals only.

Volatile Organics

In reference to question B:

At the client's request, the analytical method specified in the CAM protocol was not followed.

In reference to question H:

An MS/Dup was performed in lieu of an LCS/LCSD.

In reference to question I:

All samples were analyzed for a subset of MCP compounds per the Chain of Custody.



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

Case Narrative (continued)

Metals

L1008812-01 through -08 have elevated detection limits for Antimony and Thallium due to the dilutions required by the high concentrations of non-target analytes.

In reference to question G:

One or more of the target analytes did not achieve the requested CAM reporting limits.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

King L. Wisters Lisa Westerlind

Authorized Signature:

Title: Technical Director/Representative

Date: 06/21/10

ORGANICS



VOLATILES



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Report Date: Project Number: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: Date Collected: L1008812-01 06/11/10 08:02

Client ID: MW-3 Date Received:

06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water Analytical Method: 16,524.2 06/14/10 09:52 Analytical Date:

Analyst: TT

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--------------------------------------|----------|-----------|-------|------|-----|------------------------|
| Volatile Organics by GC/MS - Westbor | ough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| | | | | | | |



Project Name: WALPOLE PARK SOUTH **Lab Number:** L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-01 Date Collected: 06/11/10 08:02

Client ID: MW-3 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 08:02

See Narrative

06/11/10

Date Collected:

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-01

Client ID: MW-3

Sample Location: WALPOLE, MA

Parameter Result Qualifier Units RL MDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 96 | | 80-120 | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Report Date: Project Number: 12700058-003 06/21/10

SAMPLE RESULTS

Date Collected: Lab ID: L1008812-02 06/11/10 08:56

Client ID: RIZ-3

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water Analytical Method: 16,524.2 Analytical Date: 06/14/10 11:40

Analyst: TT

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbore | ough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| | | | | | | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-02 Date Collected: 06/11/10 08:56

Client ID: RIZ-3 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| VIII 022, III 1 | | | o.aop. | | Coortananto | |
|--|--------|-----------|--------|------|-------------|-----------------|
| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
| Volatile Organics by GC/MS - Westborough | ı Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 08:56

See Narrative

06/11/10

Date Collected:

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-02

Client ID: RIZ-3

Sample Location: WALPOLE, MA

Parameter Result Qualifier Units RL MDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|------------------------|------------|-----------|------------------------|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 |
| 4-Bromofluorobenzene | 96 | | 80-120 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

SAMPLE RESULTS

Lab ID: L1008812-03

Client ID: MW-2

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/14/10 12:17

Analyst: TT

Date Collected: 06/11/10 09:33 Date Received: 06/11/10

Field Prep: See Narrative

| | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westborough | Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | 1.2 | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-03 Date Collected: 06/11/10 09:33

Client ID: MW-2 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-03

Client ID: MW-2

Sample Location: WALPOLE, MA

Date Collected:

06/11/10 09:33

Date Received:

06/11/10

Field Prep:

See Narrative

Parameter Result Qualifier Units RL MDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 96 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

Report Date: 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-04

Client ID: GHC-6

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/14/10 12:54

Analyst: TT

Date Collected: 06/11/10 10:08 Date Received: 06/11/10

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|-------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbo | orough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| | | | | | | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-04 Date Collected: 06/11/10 10:08

Client ID: GHC-6 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 10:08

See Narrative

06/11/10

Project Name: Lab Number: WALPOLE PARK SOUTH L1008812

Project Number: Report Date: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-04

Client ID: GHC-6

Parameter

Sample Location: WALPOLE, MA

Field Prep: MDL Qualifier Units RL Result **Dilution Factor**

Date Collected:

Date Received:

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 96 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

SAMPLE RESULTS

Lab ID: L1008812-05

Client ID: RIZ-9

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/14/10 13:31

Analyst: TT

Date Collected: 06/11/10 11:07

Date Received: 06/11/10
Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbore | ough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-05 Date Collected: 06/11/10 11:07

Client ID: RIZ-9 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| • • • • • • • • • • • • • • • • • • • | | | | аор. | | |
|---------------------------------------|-----------|-----------|-------|------|-----|-----------------|
| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
| Volatile Organics by GC/MS - Westbo | rough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 11:07

See Narrative

06/11/10

Date Collected:

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-05

Client ID: RIZ-9

Sample Location: WALPOLE, MA

Parameter Result Qualifier Units RL MDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|------------------------|------------|-----------|------------------------|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 |
| 4-Bromofluorobenzene | 96 | | 80-120 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

Report Date: 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-06

Client ID: RIZ-10

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/14/10 14:08

Analyst: TT

Date Collected: 06/11/10 12:30

Date Received: 06/11/10
Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|-------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbo | orough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| | | | | | | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-06 Date Collected: 06/11/10 12:30

Client ID: RIZ-10 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 12:30

See Narrative

06/11/10

Date Collected:

Date Received:

Field Prep:

Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-06

Client ID: RIZ-10

Parameter

Sample Location: WALPOLE, MA

Result Qualifier Units RL MDL Dilution Factor

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|------------------------|------------|-----------|------------------------|
| 1,2-Dichlorobenzene-d4 | 103 | | 80-120 |
| 4-Bromofluorobenzene | 98 | | 80-120 |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Report Date: **Project Number:** 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: Date Collected: L1008812-07 06/11/10 13:40

Client ID: RIZ-8

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water Analytical Method: 16,524.2 06/14/10 14:46 Analytical Date:

Analyst: TT

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbore | ough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-07 Date Collected: 06/11/10 13:40

Client ID: RIZ-8 Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 13:40

Date Collected:

Project Name: Lab Number: WALPOLE PARK SOUTH L1008812

Project Number: Report Date: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-07

Client ID: RIZ-8

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter MDL Qualifier Units RL Result **Dilution Factor**

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 97 | | 80-120 | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Report Date: Project Number: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: Date Collected: L1008812-08 06/11/10 14:10

Client ID: MW-9

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Matrix: Water Analytical Method: 16,524.2 06/14/10 15:23 Analytical Date:

Analyst: TT

| Volatile Organics by GC/MS - Westborough Lab Methylene chloride ,1-Dichloroethane Chloroform Carbon tetrachloride ,2-Dichloropropane Dibromochloromethane | ND ND ND ND | ug/l ug/l | 0.50 0.50 | 1 |
|---|----------------|--------------|--------------|-------|
| ,1-Dichloroethane Chloroform Carbon tetrachloride ,2-Dichloropropane | ND ND ND | ug/l | | 1 |
| Chloroform Carbon tetrachloride ,2-Dichloropropane | ND ND | | 0.50 | |
| Carbon tetrachloride ,2-Dichloropropane | ND | | | 1 |
| ,2-Dichloropropane | | ug/l | 0.50 | 1 |
| · · | ND | ug/l | 0.50 | 1 |
| Dibromochloromethane | · ·- | ug/l | 0.50 | 1 |
| | ND | ug/l | 0.50 | 1 |
| ,1,2-Trichloroethane | ND | ug/l | 0.50 | 1 |
| etrachloroethene | ND | ug/l | 0.50 | 1 |
| Chlorobenzene | ND | ug/l | 0.50 | 1 |
| richlorofluoromethane | ND | ug/l | 0.50 | 1 |
| ,2-Dichloroethane | ND | ug/l | 0.50 | 1 |
| ,1,1-Trichloroethane | ND | ug/l | 0.50 | 1 |
| Bromodichloromethane | ND | ug/l | 0.50 | 1 |
| rans-1,3-Dichloropropene | ND | ug/l | 0.50 | 1 |
| sis-1,3-Dichloropropene | ND | ug/l | 0.50 | 1 |
| Bromoform | ND | ug/l | 0.50 | 1 |
| ,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 1 |
| Benzene | ND | ug/l | 0.50 | 1 |
| oluene | ND | ug/l | 0.50 | 1 |
| Ethylbenzene | ND | ug/l | 0.50 | 1 |
| /m-Xylene | ND | ug/l | 0.50 | 1 |
| Chloromethane | ND | ug/l | 0.50 | 1 |
| Bromomethane | ND | ug/l | 0.50 | 1 |
| /inyl chloride | ND | ug/l | 0.50 | 1 |
| Chloroethane | ND | ug/l | 0.50 | 1 |
| ,1-Dichloroethene | ND | ug/l | 0.50 | 1 |
| rans-1,2-Dichloroethene | ND | ug/l | 0.50 | 1 |
| sis-1,2-Dichloroethene | ND | ug/l | 0.50 | 1 |
| richloroethene | ND | ug/l | 0.50 | 1 |
| ,2-Dichlorobenzene | ND | ug/l | 0.50 | 1 |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-08 Date Collected: 06/11/10 14:10

Client ID: Date Received: 06/11/10

Sample Location: WALPOLE, MA Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|------------------------------------|------------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westb | orough Lab | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Styrene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 14:10

Date Collected:

Project Name: Lab Number: WALPOLE PARK SOUTH L1008812

Project Number: Report Date: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-08

Client ID: MW-9

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: See Narrative

Parameter MDL Qualifier Units RL Result **Dilution Factor**

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 104 | | 80-120 | |
| 4-Bromofluorobenzene | 97 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-09

Client ID: 20100611-TRIP BLANK

Sample Location: WALPOLE, MA

Matrix: Water
Analytical Method: 16,524.2
Analytical Date: 06/14/10 16:00

Analyst: TT

| Date Collected: | 06/11/10 00:00 |
|-----------------|----------------|
| Date Received: | 06/11/10 |
| Field Prep: | Not Specified |
| | |
| | |
| | |
| | |

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------------|----------|-----------|-------|------|-----|-----------------|
| Volatile Organics by GC/MS - Westbore | ough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| Chloroform | ND | | ug/l | 0.50 | | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | | 1 |
| Chlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| Bromoform | ND | | ug/l | 0.50 | | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| Benzene | ND | | ug/l | 0.50 | | 1 |
| Toluene | ND | | ug/l | 0.50 | | 1 |
| Ethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p/m-Xylene | ND | | ug/l | 0.50 | | 1 |
| Chloromethane | ND | | ug/l | 0.50 | | 1 |
| Bromomethane | ND | | ug/l | 0.50 | | 1 |
| Vinyl chloride | ND | | ug/l | 0.50 | | 1 |
| Chloroethane | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: Report Date: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: Date Collected: L1008812-09 06/11/10 00:00

Client ID: 20100611-TRIP BLANK Date Received: 06/11/10 Field Prep: Not Specified

Sample Location: WALPOLE, MA

| 1,3-Dichlorobenzene ND ug/l 0.50 1 1,4-Dichlorobenzene ND ug/l 0.50 1 1,4-Dichlorobenzene ND ug/l 0.50 1 1,5-Dichlorobenzene ND ug/l 0.50 1 0-Xylene ND ug/l 0.50 1 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloroethane ND ug/l 0.50 1 1,1,1,1,1,1,1,1,1,1,1,1,1,1,1,1,1,1,1, | Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|-------------------------------------|------------|-----------|-------|------|-----|-----------------|
| 1,4-Dichlorobenzene ND ugl 0.50 1 Styrene ND ugl 0.50 1 c-Xylene ND ugl 0.50 1 1,1-Dichloropropene ND ugl 0.50 1 1,1-1,2-Tetrachloropethane ND ugl 0.50 1 1,2,3-Trichloropropane ND ugl 0.50 1 Bromochloromethane ND ugl 0.50 1 Bromochloromethane ND ugl 0.50 1 I-Butylbenzene ND ugl 0.50 1 Dichlorodifluoromethane ND ugl 0.50 1 Hexachlorobutadiene ND ugl 0.50 1 Isopropylbenzene ND ugl 0.50 1 Isopropylbenzene ND ugl 0.50 1 | Volatile Organics by GC/MS - Westbo | orough Lab | | | | | |
| Skyrene ND ugl 0.50 1 c-Xylene ND ugl 0.50 1 1.1-Dichloropropene ND ugl 0.50 1 2,2-Dichloropropane ND ugl 0.50 1 1,1,1,2-Etarchloroethane ND ugl 0.50 1 1,1,2-Etrachloropropane ND ugl 0.50 1 1,2,2-Trichloropropane ND ugl 0.50 1 Bromochloromethane ND ugl 0.50 1 Bromochloromethane ND ugl 0.50 1 Hexachlorobutadiene ND ugl 0.50 1 Hexachlorobutadiene ND ugl 0.50 1 p-Isopropylibulene ND ugl 0.50 1 Naphthalene ND ugl 0.50 1 | 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| o-Xylene ND ug/l 0.50 1 1.1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1.1,1,2-Tetrachloropethane ND ug/l 0.50 1 1.2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Dichlorodifloromethane ND ug/l 0.50 1 Dichlorodifloromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 <td>1,4-Dichlorobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td></td> <td>1</td> | 1,4-Dichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,1-Dichloropropene ND ug/l 0.50 1 2,2-Dichloropropane ND ug/l 0.50 1 1,1,1,2-Tetrachloropethane ND ug/l 0.50 1 1,2,3-Trichloropropane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 Bromochloromethane ND ug/l 0.50 1 D-Butylbenzene ND ug/l 0.50 1 D-Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 see-Butylbenzene ND ug/l 0.50 1 | Styrene | ND | | ug/l | 0.50 | | 1 |
| 2,2-Dichloropropane ND ug/l 0,50 1 1,1,1,2-Tetrachloroethane ND ug/l 0,50 1 1,2,3-Trichloropropane ND ug/l 0,50 1 Bromochloromethane ND ug/l 0,50 1 n-Butybenzene ND ug/l 0,50 1 Dichlorodiffuoromethane ND ug/l 0,50 1 Hexachlorobutadiene ND ug/l 0,50 1 Bopropyblenzene ND ug/l 0,50 1 Isopropyblenzene ND ug/l 0,50 1 Naphthalene ND ug/l 0,50 1 n-Propyblenzene ND ug/l 0,50 1 sec-Butybenzene ND ug/l 0,50 1 tert-Butybenzene ND ug/l 0,50 1 <td>o-Xylene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td></td> <td>1</td> | o-Xylene | ND | | ug/l | 0.50 | | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 0,50 1 1,2,3-Trichloropropane ND ug/l 0,50 1 Bromochloromethane ND ug/l 0,50 1 n-Butylbenzene ND ug/l 0,50 1 Dichlorodiffluoromethane ND ug/l 0,50 1 Hexachlorobutadiene ND ug/l 0,50 1 Isopropylbenzene ND ug/l 0,50 1 Isopropylbenzene ND ug/l 0,50 1 Isopropylbenzene ND ug/l 0,50 1 Naphthalene ND ug/l 0,50 1 n-Propylbenzene ND ug/l 0,50 1 sec-Butylbenzene ND ug/l 0,50 1 tert-Butylbenzene ND ug/l 0,50 1 < | 1,1-Dichloropropene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichloropropane ND ug/l 0.50 1 | 2,2-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| Bromochloromethane ND ug/l 0.50 1 n-Butylbenzene ND ug/l 0.50 1 Dichlorodiffluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 N-Propylbenzene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.50 | | 1 |
| n-Butylbenzene ND ug/l 0.50 1 Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 | 1,2,3-Trichloropropane | ND | | ug/l | 0.50 | | 1 |
| Dichlorodifluoromethane ND ug/l 0.50 1 Hexachlorobutadiene ND ug/l 0.50 1 Isopropylbenzene ND ug/l 0.50 1 p-Isopropyltoluene ND ug/l 0.50 1 Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 </td <td>Bromochloromethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td></td> <td>1</td> | Bromochloromethane | ND | | ug/l | 0.50 | | 1 |
| Hexachlorobutadiene ND Ug/l 0.50 1 | n-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| Sopropyltoluene ND | Dichlorodifluoromethane | ND | | ug/l | 0.50 | | 1 |
| P-Isopropyltoluene ND ug/l 0.50 1 | Hexachlorobutadiene | ND | | ug/l | 0.50 | | 1 |
| Naphthalene ND ug/l 0.50 1 n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 | Isopropylbenzene | ND | | ug/l | 0.50 | | 1 |
| n-Propylbenzene ND ug/l 0.50 1 sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 </td <td>p-Isopropyltoluene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td></td> <td>1</td> | p-Isopropyltoluene | ND | | ug/l | 0.50 | | 1 |
| sec-Butylbenzene ND ug/l 0.50 1 tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Naphthalene | ND | | ug/l | 0.50 | | 1 |
| tert-Butylbenzene ND ug/l 0.50 1 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 0-Chlorotoluene ND ug/l 0.50 1 1 p-Chlorotoluene ND ug/l 0.50 1 1 p-Chlorotoluene ND ug/l 0.50 1 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | n-Propylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,3-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | sec-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trichlorobenzene ND ug/l 0.50 1 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | tert-Butylbenzene | ND | | ug/l | 0.50 | | 1 |
| 1,2,4-Trimethylbenzene ND ug/l 0.50 1 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,3-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| 1,3,5-Trimethylbenzene ND ug/l 0.50 1 Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trichlorobenzene | ND | | ug/l | 0.50 | | 1 |
| Bromobenzene ND ug/l 0.50 1 o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2,4-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| o-Chlorotoluene ND ug/l 0.50 1 p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | 1,3,5-Trimethylbenzene | ND | | ug/l | 0.50 | | 1 |
| p-Chlorotoluene ND ug/l 0.50 1 Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Bromobenzene | ND | | ug/l | 0.50 | | 1 |
| Dibromomethane ND ug/l 0.50 1 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | o-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromoethane ND ug/l 0.50 1 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | p-Chlorotoluene | ND | | ug/l | 0.50 | | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 0.50 1 1,3-Dichloropropane ND ug/l 0.50 1 | Dibromomethane | ND | | ug/l | 0.50 | | 1 |
| 1,3-Dichloropropane ND ug/l 0.50 1 | 1,2-Dibromoethane | ND | | ug/l | 0.50 | | 1 |
| | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.50 | | 1 |
| Methyl tert butyl ether ND ug/l 0.50 1 | 1,3-Dichloropropane | ND | | ug/l | 0.50 | | 1 |
| | Methyl tert butyl ether | ND | | ug/l | 0.50 | | 1 |

| Tentatively Identified Compounds | | | |
|-------------------------------------|----|------|---|
| No Tentatively Identified Compounds | ND | ug/l | 1 |



06/11/10 00:00

Date Collected:

Project Name: Lab Number: WALPOLE PARK SOUTH L1008812

Project Number: Report Date: 12700058-003 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-09

Client ID: 20100611-TRIP BLANK

Date Received: 06/11/10 Sample Location: WALPOLE, MA Field Prep: Not Specified

Parameter MDL Result Qualifier Units RL **Dilution Factor**

Volatile Organics by GC/MS - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|------------------------|------------|-----------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 100 | | 80-120 | |
| 4-Bromofluorobenzene | 97 | | 80-120 | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/14/10 07:26

Analyst: TT

| arameter | Result | Qualifier | Units | RL | MDL |
|---------------------------|-------------------|------------------|-------|------------|---------|
| olatile Organics by GC/MS | - Westborough Lal | o for sample(s): | 01-09 | Batch: WG4 | 17697-2 |
| Methylene chloride | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethane | ND | | ug/l | 0.50 | |
| Chloroform | ND | | ug/l | 0.50 | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | |
| 1,2-Dichloropropane | ND | | ug/l | 0.50 | |
| Dibromochloromethane | ND | | ug/l | 0.50 | |
| 1,1,2-Trichloroethane | ND | | ug/l | 0.50 | |
| Tetrachloroethene | ND | | ug/l | 0.50 | |
| Chlorobenzene | ND | | ug/l | 0.50 | |
| Trichlorofluoromethane | ND | | ug/l | 0.50 | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | |
| 1,1,1-Trichloroethane | ND | | ug/l | 0.50 | |
| Bromodichloromethane | ND | | ug/l | 0.50 | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | |
| Bromoform | ND | | ug/l | 0.50 | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | |
| Benzene | ND | | ug/l | 0.50 | |
| Toluene | ND | | ug/l | 0.50 | |
| Ethylbenzene | ND | | ug/l | 0.50 | |
| p/m-Xylene | ND | | ug/l | 0.50 | |
| Chloromethane | ND | | ug/l | 0.50 | |
| Bromomethane | ND | | ug/l | 0.50 | |
| Vinyl chloride | ND | | ug/l | 0.50 | |
| Chloroethane | ND | | ug/l | 0.50 | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | |
| trans-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| cis-1,2-Dichloroethene | ND | | ug/l | 0.50 | |
| Trichloroethene | ND | | ug/l | 0.50 | |
| 1,2-Dichlorobenzene | ND | | ug/l | 0.50 | |
| 1,3-Dichlorobenzene | ND | | ug/l | 0.50 | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

Method Blank Analysis Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/14/10 07:26

Analyst: TT

| Parameter | Result | Qualifier | Units | R | L MDL |
|-----------------------------|-------------------|----------------|-------|--------|------------|
| olatile Organics by GC/MS | - Westborough Lab | for sample(s): | 01-09 | Batch: | WG417697-2 |
| 1,4-Dichlorobenzene | ND | | ug/l | 0.5 | 50 |
| Styrene | ND | | ug/l | 0.5 | 50 |
| o-Xylene | ND | | ug/l | 0.5 | 50 |
| 1,1-Dichloropropene | ND | | ug/l | 0.5 | 50 |
| 2,2-Dichloropropane | ND | | ug/l | 0.5 | 50 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 0.5 | 50 |
| 1,2,3-Trichloropropane | ND | | ug/l | 0.5 | 50 |
| Bromochloromethane | ND | | ug/l | 0.5 | 50 |
| n-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| Dichlorodifluoromethane | ND | | ug/l | 0.5 | 50 |
| Hexachlorobutadiene | ND | | ug/l | 0.5 | 50 |
| Isopropylbenzene | ND | | ug/l | 0.5 | 50 |
| p-Isopropyltoluene | ND | | ug/l | 0.5 | 50 |
| Naphthalene | ND | | ug/l | 0.5 | 50 |
| n-Propylbenzene | ND | | ug/l | 0.5 | 50 |
| sec-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| tert-Butylbenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 0.5 | 50 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 0.5 | 50 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 0.5 | 50 |
| Bromobenzene | ND | | ug/l | 0.5 | 50 |
| o-Chlorotoluene | ND | | ug/l | 0.5 | 50 |
| p-Chlorotoluene | ND | | ug/l | 0.5 | 50 |
| Dibromomethane | ND | | ug/l | 0.5 | 50 |
| 1,2-Dibromoethane | ND | | ug/l | 0.5 | 50 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 0.5 | 50 |
| 1,3-Dichloropropane | ND | | ug/l | 0.5 | 50 |
| Methyl tert butyl ether | ND | | ug/l | 0.5 | 50 |
| | | | | | |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

Method Blank Analysis
Batch Quality Control

Analytical Method: 16,524.2 Analytical Date: 06/14/10 07:26

Analyst: TT

| Parameter | Result | Qualifier | Units | RL | MDL | |
|----------------------------|----------------|--------------|------------|------------|---------|--|
| Valatila Organica by CC/MS | Weatherough Le | h for comple | (a): 01 00 | Potob: WC4 | 17607 2 | |

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-09 Batch: WG417697-2

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l

| | | 4 | Acceptance | |
|------------------------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | Criteria | |
| | | | | |
| 1,2-Dichlorobenzene-d4 | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 97 | | 80-120 | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | LCS %Recovery | Qual | | CSD covery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|------------|-------|---------------|-----------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough I | _ab Associated | sample(s): | 01-09 | Batch: | WG417697- | 1 | | | |
| Methylene chloride | 107 | | | - | | 70-130 | - | | |
| 1,1-Dichloroethane | 106 | | | - | | 70-130 | - | | |
| Chloroform | 102 | | | - | | 70-130 | - | | |
| Carbon tetrachloride | 86 | | | - | | 70-130 | - | | |
| 1,2-Dichloropropane | 107 | | | - | | 70-130 | - | | |
| Dibromochloromethane | 95 | | | - | | 70-130 | - | | |
| 1,1,2-Trichloroethane | 102 | | | - | | 70-130 | - | | |
| Tetrachloroethene | 106 | | | - | | 70-130 | - | | |
| Chlorobenzene | 97 | | | - | | 70-130 | - | | |
| Trichlorofluoromethane | 102 | | | - | | 70-130 | - | | |
| 1,2-Dichloroethane | 99 | | | - | | 70-130 | - | | |
| 1,1,1-Trichloroethane | 99 | | | - | | 70-130 | - | | |
| Bromodichloromethane | 94 | | | - | | 70-130 | - | | |
| trans-1,3-Dichloropropene | 79 | | | - | | 70-130 | - | | |
| cis-1,3-Dichloropropene | 84 | | | - | | 70-130 | - | | |
| Bromoform | 86 | | | - | | 70-130 | - | | |
| 1,1,2,2-Tetrachloroethane | 94 | | | - | | 70-130 | - | | |
| Benzene | 109 | | | - | | 70-130 | - | | |
| Toluene | 108 | | | - | | 70-130 | - | | |
| Ethylbenzene | 98 | | | - | | 70-130 | - | | |
| p/m-Xylene | 100 | | | - | | 70-130 | - | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | LCS %Recovery (| LCS Qual %Reco | | %Recovery Limits | RPD | Qual | RPD Limits |
|---|--------------------|-------------------|---------------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough L | ab Associated sar | mple(s): 01-09 l | Batch: WG4176 | 697-1 | | | |
| Chloromethane | 119 | - | | 70-130 | - | | |
| Bromomethane | 116 | - | | 70-130 | - | | |
| Vinyl chloride | 117 | - | | 70-130 | - | | |
| Chloroethane | 112 | - | | 70-130 | - | | |
| 1,1-Dichloroethene | 109 | - | | 70-130 | - | | |
| trans-1,2-Dichloroethene | 109 | - | | 70-130 | - | | |
| cis-1,2-Dichloroethene | 104 | - | | 70-130 | - | | |
| Trichloroethene | 103 | - | | 70-130 | - | | |
| 1,2-Dichlorobenzene | 91 | - | | 70-130 | - | | |
| 1,3-Dichlorobenzene | 94 | - | | 70-130 | - | | |
| 1,4-Dichlorobenzene | 93 | - | | 70-130 | - | | |
| Styrene | 97 | - | | 70-130 | - | | |
| o-Xylene | 97 | - | | 70-130 | - | | |
| 1,1-Dichloropropene | 104 | - | | 70-130 | - | | |
| 2,2-Dichloropropane | 78 | - | | 70-130 | - | | |
| 1,1,1,2-Tetrachloroethane | 92 | - | | 70-130 | - | | |
| 1,2,3-Trichloropropane | 92 | - | | 70-130 | - | | |
| Bromochloromethane | 101 | - | | 70-130 | - | | |
| n-Butylbenzene | 96 | - | | 70-130 | - | | |
| Dichlorodifluoromethane | 122 | - | | 70-130 | - | | |
| Hexachlorobutadiene | 98 | - | | 70-130 | - | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | LCS %Recovery | Qual | | CSD covery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|---|------------------|------------|-------|---------------|-----------|---------------------|-----|------|------------|
| olatile Organics by GC/MS - Westborough L | ab Associated | sample(s): | 01-09 | Batch: | WG417697- | -1 | | | |
| Isopropylbenzene | 97 | | | - | | 70-130 | - | | |
| p-Isopropyltoluene | 98 | | | - | | 70-130 | - | | |
| Naphthalene | 74 | | | - | | 70-130 | - | | |
| n-Propylbenzene | 98 | | | - | | 70-130 | - | | |
| sec-Butylbenzene | 97 | | | - | | 70-130 | - | | |
| tert-Butylbenzene | 97 | | | - | | 70-130 | - | | |
| 1,2,3-Trichlorobenzene | 82 | | | - | | 70-130 | - | | |
| 1,2,4-Trichlorobenzene | 86 | | | - | | 70-130 | - | | |
| 1,2,4-Trimethylbenzene | 98 | | | - | | 70-130 | - | | |
| 1,3,5-Trimethylbenzene | 98 | | | - | | 70-130 | - | | |
| Bromobenzene | 96 | | | - | | 70-130 | - | | |
| o-Chlorotoluene | 100 | | | - | | 70-130 | - | | |
| p-Chlorotoluene | 99 | | | - | | 70-130 | - | | |
| Dibromomethane | 98 | | | - | | 70-130 | - | | |
| 1,2-Dibromoethane | 91 | | | - | | 70-130 | - | | |
| 1,2-Dibromo-3-chloropropane | 78 | | | - | | 70-130 | - | | |
| 1,3-Dichloropropane | 100 | | | - | | 70-130 | - | | |
| Methyl tert butyl ether | 101 | | | - | | 70-130 | - | | |



Lab Control Sample Analysis

WALPOLE PARK SOUTH

Batch Quality Control

Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

LCS LCSD %Recovery

Parameter %Recovery Qual %Recovery Qual Limits RPD Qual RPD Limits

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-09 Batch: WG417697-1

| Surrogate | LCS %Recovery | Qual | LCSD %Recovery | Qual | Acceptance Criteria | |
|------------------------|------------------|------|-------------------|------|------------------------|--|
| 1,2-Dichlorobenzene-d4 | 98 | | | | 80-120 | |
| 4-Bromofluorobenzene | 98 | | | | 80-120 | |



Project Name:

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| rameter | Native Sample | MS Added | MS Found | MS %Recovery | MSD Qual Found | MSD %Recovery | Recovery Qual Limits | RPD | | RPD <u>Limit</u> s |
|--------------------------------------|------------------|-------------|-------------|-----------------|--------------------|------------------|-------------------------|------|------------|-----------------------|
| olatile Organics by GC/MS - ample | - Westborough | Lab Associ | ated sample | (s): 01-09 Q | QC Batch ID: WG417 | 7697-3 QC | Sample: L1008822 | 2-01 | Client ID: | MS |
| Methylene chloride | ND | 4 | 4.5 | 112 | | - | 70-130 | - | | 20 |
| 1,1-Dichloroethane | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | | 20 |
| Chloroform | 27 | 4 | 31 | 102 | - | - | 70-130 | - | | 20 |
| Carbon tetrachloride | ND | 4 | 4.0 | 101 | - | - | 70-130 | - | | 20 |
| 1,2-Dichloropropane | ND | 4 | 4.5 | 113 | - | - | 70-130 | - | | 20 |
| Dibromochloromethane | 4.8 | 4 | 8.8 | 100 | - | - | 70-130 | - | | 20 |
| 1,1,2-Trichloroethane | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | | 20 |
| Tetrachloroethene | ND | 4 | 4.6 | 116 | - | - | 70-130 | - | | 20 |
| Chlorobenzene | ND | 4 | 4.1 | 104 | - | - | 70-130 | - | | 20 |
| Trichlorofluoromethane | ND | 4 | 4.6 | 114 | - | - | 70-130 | - | | 20 |
| 1,2-Dichloroethane | ND | 4 | 4.1 | 102 | - | - | 70-130 | - | | 20 |
| 1,1,1-Trichloroethane | ND | 4 | 4.3 | 108 | - | - | 70-130 | - | | 20 |
| Bromodichloromethane | 10 | 4 | 14 | 100 | - | - | 70-130 | - | | 20 |
| trans-1,3-Dichloropropene | ND | 4 | 3.1 | 77 | - | - | 70-130 | - | | 20 |
| cis-1,3-Dichloropropene | ND | 4 | 4.0 | 99 | - | - | 70-130 | - | | 20 |
| Bromoform | 0.59 | 4 | 4.2 | 91 | | - | 70-130 | - | | 20 |
| 1,1,2,2-Tetrachloroethane | ND | 4 | 3.8 | 96 | | - | 70-130 | - | | 20 |
| Benzene | ND | 4 | 4.5 | 112 | | - | 70-130 | - | | 20 |
| Toluene | ND | 4 | 4.5 | 112 | | - | 70-130 | - | | 20 |
| Ethylbenzene | ND | 4 | 4.4 | 110 | | - | 70-130 | - | | 20 |
| p/m-Xylene | 1.2 | 8 | 9.4 | 102 | | - | 70-130 | - | | 20 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | | Recovery Limits | RPD | Qual | RPD Limits |
|--------------------------------------|------------------|-------------|--------------|-----------------|---------|--------------|------------------|--------|--------------------|------|------------|---------------|
| Volatile Organics by GC/MS Sample | - Westborough | Lab Assoc | iated sample | (s): 01-09 Q | C Batch | ID: WG41 | 7697-3 QC | Sample | : L100882 | 2-01 | Client ID: | MS |
| Chloromethane | ND | 4 | 4.0 | 101 | | - | - | | 70-130 | - | | 20 |
| Bromomethane | ND | 4 | 4.6 | 116 | | - | - | | 70-130 | - | | 20 |
| Vinyl chloride | ND | 4 | 5.9 | 148 | Q | - | - | | 70-130 | - | | 20 |
| Chloroethane | ND | 4 | 4.9 | 122 | | - | - | | 70-130 | - | | 20 |
| 1,1-Dichloroethene | ND | 4 | 4.6 | 115 | | - | - | | 70-130 | - | | 20 |
| trans-1,2-Dichloroethene | ND | 4 | 4.5 | 113 | | - | - | | 70-130 | - | | 20 |
| cis-1,2-Dichloroethene | ND | 4 | 4.6 | 116 | | - | - | | 70-130 | - | | 20 |
| Trichloroethene | ND | 4 | 4.3 | 108 | | - | - | | 70-130 | - | | 20 |
| 1,2-Dichlorobenzene | ND | 4 | 3.8 | 95 | | - | - | | 70-130 | - | | 20 |
| 1,3-Dichlorobenzene | ND | 4 | 3.8 | 95 | | - | - | | 70-130 | - | | 20 |
| 1,4-Dichlorobenzene | ND | 4 | 3.7 | 94 | | - | - | | 70-130 | - | | 20 |
| Styrene | ND | 4 | 3.7 | 93 | | - | - | | 70-130 | - | | 20 |
| o-Xylene | ND | 4 | 4.4 | 110 | | - | - | | 70-130 | - | | 20 |
| 1,1-Dichloropropene | ND | 4 | 4.6 | 116 | | - | - | | 70-130 | - | | 20 |
| 2,2-Dichloropropane | ND | 4 | 3.7 | 94 | | - | - | | 70-130 | - | | 20 |
| 1,1,1,2-Tetrachloroethane | ND | 4 | 3.9 | 98 | | - | - | | 70-130 | - | | 20 |
| 1,2,3-Trichloropropane | ND | 4 | 3.6 | 91 | | - | - | | 70-130 | - | | 20 |
| Bromochloromethane | ND | 4 | 4.5 | 112 | | - | - | | 70-130 | - | | 20 |
| n-Butylbenzene | ND | 4 | 3.9 | 98 | | - | - | | 70-130 | - | | 20 |
| Dichlorodifluoromethane | ND | 4 | 4.2 | 105 | | - | - | | 70-130 | - | | 20 |
| Hexachlorobutadiene | ND | 4 | 3.9 | 99 | | - | - | | 70-130 | - | | 20 |
| | | | | | | | | | | | | |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | Native Sample | MS Added | MS Found | MS %Recovery | Qual | MSD Found | MSD %Recovery | Recovery Qual Limits | RPD | Qual | RPD <u>Limit</u> s |
|-----------------------------|------------------|-------------|-------------|-----------------|------------|--------------|------------------|-------------------------|------|------------|-----------------------|
| olatile Organics by GC/MS - | Westborough | Lab Associ | ated sample | (s): 01-09 Q | C Batch II | D: WG417 | 7697-3 QC | Sample: L100882 | 2-01 | Client ID: | MS |
| Isopropylbenzene | ND | 4 | 3.6 | 90 | | - | - | 70-130 | - | | 20 |
| p-Isopropyltoluene | ND | 4 | 3.9 | 98 | | - | - | 70-130 | - | | 20 |
| Naphthalene | 0.59 | 4 | 3.9 | 82 | | - | - | 70-130 | - | | 20 |
| n-Propylbenzene | ND | 4 | 4.1 | 102 | | - | - | 70-130 | - | | 20 |
| sec-Butylbenzene | ND | 4 | 4.1 | 102 | | - | - | 70-130 | - | | 20 |
| tert-Butylbenzene | ND | 4 | 4.1 | 102 | | - | - | 70-130 | - | | 20 |
| 1,2,3-Trichlorobenzene | ND | 4 | 3.4 | 86 | | - | - | 70-130 | - | | 20 |
| 1,2,4-Trichlorobenzene | ND | 4 | 3.4 | 85 | | - | - | 70-130 | - | | 20 |
| 1,2,4-Trimethylbenzene | ND | 4 | 4.0 | 101 | | - | - | 70-130 | - | | 20 |
| 1,3,5-Trimethylbenzene | ND | 4 | 4.0 | 100 | | - | - | 70-130 | - | | 20 |
| Bromobenzene | ND | 4 | 3.9 | 98 | | - | - | 70-130 | - | | 20 |
| o-Chlorotoluene | ND | 4 | 4.2 | 104 | | - | - | 70-130 | - | | 20 |
| p-Chlorotoluene | ND | 4 | 3.9 | 98 | | - | - | 70-130 | - | | 20 |
| Dibromomethane | ND | 4 | 4.3 | 108 | | - | - | 70-130 | - | | 20 |
| 1,2-Dibromoethane | ND | 4 | 3.7 | 93 | | - | - | 70-130 | - | | 20 |
| 1,2-Dibromo-3-chloropropane | ND | 4 | 3.4 | 86 | | - | - | 70-130 | - | | 20 |
| 1,3-Dichloropropane | ND | 4 | 4.1 | 103 | | - | - | 70-130 | - | | 20 |
| Methyl tert butyl ether | ND | 4 | 4.2 | 104 | | - | - | 70-130 | - | | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

| | Native | MS | MS | MS | | MSD | MSD | | Recovery | | | RPD |
|-----------|--------|-------|-------|-----------|------|-------|-----------|------|----------|-----|------|--------|
| Parameter | Sample | Added | Found | %Recovery | Qual | Found | %Recovery | Qual | Limits | RPD | Qual | Limits |

Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-09 QC Batch ID: WG417697-3 QC Sample: L1008822-01 Client ID: MS Sample

| | MS | MSD | Acceptance |
|------------------------|----------------------|----------------------|------------|
| Surrogate | % Recovery Qualifier | % Recovery Qualifier | Criteria |
| 1,2-Dichlorobenzene-d4 | 100 | | 80-120 |
| 4-Bromofluorobenzene | 99 | | 80-120 |

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

L1008812 06/21/10 Report Date:

Lab Number:

| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-09 QC Batch ID: WG417697-4 QC Sample: L1008812-01 Client ID: MW-3 Methylene chloride ND ND ND ugil NC 20 1,1-Dichloroethane ND ND ND ugil NC 20 Carbon tetrachloride ND ND ND ugil NC 20 1,2-Dichloropropane ND ND ND ugil NC 20 Dibromochloromethane ND ND ND ugil NC 20 1,1,2-Trichloroethane ND ND ND ugil NC 20 Chlorobenzene ND ND ND ugil NC 20 1,2-Dichloroethane ND ND ND ugil NC 20 1,1,1-Trichloroethane ND ND ND ugil NC 20 Bromodichloromethane ND ND ND ugil | arameter | Native Sample | Duplicate Sample | Units | RPD | Qual RPD Limits |
|---|---|-----------------------------|------------------------|-------------|-----------|------------------------|
| 1,1-Dichloroethane | olatile Organics by GC/MS - Westborough Lab | Associated sample(s): 01-09 | QC Batch ID: WG417697- | 4 QC Sample | e: L10088 | 312-01 Client ID: MW-3 |
| Chloroform ND ND ug/l NC 20 Carbon tetrachloride ND ND ND ug/l NC 20 1,2-Dichloropropane ND ND ND ug/l NC 20 Dibromochloromethane ND ND ND ug/l NC 20 1,1,2-Trichloroethane ND ND ND ug/l NC 20 Tetrachloroethane ND ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND <td< td=""><td>Methylene chloride</td><td>ND</td><td>ND</td><td>ug/l</td><td>NC</td><td>20</td></td<> | Methylene chloride | ND | ND | ug/l | NC | 20 |
| Carbon tetrachloride ND ND ug/l NC 20 1,2-Dichloropropane ND ND ND ug/l NC 20 Dibromochloromethane ND ND ND ug/l NC 20 1,1,2-Trichloroethane ND ND ND ug/l NC 20 Tetrachloroethane ND ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 Benzene ND ND ND ug | 1,1-Dichloroethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloropropane ND ND ug/l NC 20 Dibromochloromethane ND ND ND ug/l NC 20 1,1,2-Trichloroethane ND ND ND ug/l NC 20 Tetrachloroethane ND ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ND ug/l NC 20 1,2-Dichloroethane ND ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND <td>Chloroform</td> <td>ND</td> <td>ND</td> <td>ug/l</td> <td>NC</td> <td>20</td> | Chloroform | ND | ND | ug/l | NC | 20 |
| Dibromochloromethane ND ND ug/l NC 20 1,1,2-Trichloroethane ND ND ND ug/l NC 20 Tetrachloroethene ND ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ND ug/l NC 20 1,2-Dichloroethane ND ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND <t< td=""><td>Carbon tetrachloride</td><td>ND</td><td>ND</td><td>ug/l</td><td>NC</td><td>20</td></t<> | Carbon tetrachloride | ND | ND | ug/l | NC | 20 |
| 1,1,2-Trichloroethane ND ND ug/l NC 20 Tetrachloroethene ND ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ND ug/l NC 20 1,2-Dichloroethane ND ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | 1,2-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Tetrachloroethene ND ND ug/l NC 20 Chlorobenzene ND ND ND ug/l NC 20 Trichlorofluoromethane ND ND ug/l NC 20 1,2-Dichloroethane ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | Dibromochloromethane | ND | ND | ug/l | NC | 20 |
| Chlorobenzene ND ND ug/l NC 20 Trichlorofluoromethane ND ND ug/l NC 20 1,2-Dichloroethane ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 cis-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | 1,1,2-Trichloroethane | ND | ND | ug/l | NC | 20 |
| Trichlorofluoromethane ND ND ug/l NC 20 1,2-Dichloroethane ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ug/l NC 20 Bromodichloromethane ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | Tetrachloroethene | ND | ND | ug/l | NC | 20 |
| 1,2-Dichloroethane ND ND ug/l NC 20 1,1,1-Trichloroethane ND ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | Chlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,1,1-Trichloroethane ND ND ug/l NC 20 Bromodichloromethane ND ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 cis-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | Trichlorofluoromethane | ND | ND | ug/l | NC | 20 |
| Bromodichloromethane ND ND ug/l NC 20 trans-1,3-Dichloropropene ND ND ND ug/l NC 20 cis-1,3-Dichloropropene ND ND ND ug/l NC 20 Bromoform ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | 1,2-Dichloroethane | ND | ND | ug/l | NC | 20 |
| trans-1,3-Dichloropropene ND ND ug/l NC 20 cis-1,3-Dichloropropene ND ND ug/l NC 20 Bromoform ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ug/l NC 20 Benzene ND ND ND ug/l NC 20 | 1,1,1-Trichloroethane | ND | ND | ug/l | NC | 20 |
| cis-1,3-Dichloropropene ND ND ug/l NC 20 Bromoform ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ug/l NC 20 Benzene ND ND ug/l NC 20 | Bromodichloromethane | ND | ND | ug/l | NC | 20 |
| Bromoform ND ND ug/l NC 20 1,1,2,2-Tetrachloroethane ND ND ug/l NC 20 Benzene ND ND ug/l NC 20 | trans-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| 1,1,2,2-Tetrachloroethane ND ND ug/l NC 20 Benzene ND ND ug/l NC 20 | cis-1,3-Dichloropropene | ND | ND | ug/l | NC | 20 |
| Benzene ND ND ug/l NC 20 | Bromoform | ND | ND | ug/l | NC | 20 |
| | 1,1,2,2-Tetrachloroethane | ND | ND | ug/l | NC | 20 |
| Toluene ND ND ug/l NC 20 | Benzene | ND | ND | ug/l | NC | 20 |
| | Toluene | ND | ND | ug/l | NC | 20 |



ND

ug/l

NC

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812 **Report Date:** 06/21/10

Native Sample Duplicate Sample Units RPD RPD Limits Parameter Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-09 QC Batch ID: WG417697-4 QC Sample: L1008812-01 Client ID: MW-3 NC Ethylbenzene ND ND ug/l 20 p/m-Xylene ND ND ug/l NC 20 Chloromethane ND ND ug/l NC 20 Bromomethane ND ND ug/l NC 20 Vinyl chloride ND ND ug/l NC 20 Chloroethane ND ND ug/l NC 20 1,1-Dichloroethene ND ND ug/l NC 20 trans-1,2-Dichloroethene ND ND ug/l NC 20 cis-1,2-Dichloroethene ND ND ug/l NC 20 Trichloroethene ND ND ug/l NC 20 1,2-Dichlorobenzene ND ND ug/l NC 20 1.3-Dichlorobenzene ND ND ug/l NC 20 1,4-Dichlorobenzene ND ND ug/l NC 20 Styrene ND ND ug/l NC 20 o-Xylene ND ND ug/l NC 20 1,1-Dichloropropene ND ND ug/l NC 20 2,2-Dichloropropane ND ND ug/l NC 20 1,1,1,2-Tetrachloroethane ND ND ug/l NC 20

ND



20

1,2,3-Trichloropropane

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

| arameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|---|-----------------------------|----------------------|-------------|----------------|-------------------|
| olatile Organics by GC/MS - Westborough Lab | Associated sample(s): 01-09 | QC Batch ID: WG41769 | 7-4 QC Samp | le: L1008812-0 | 1 Client ID: MW-3 |
| Bromochloromethane | ND | ND | ug/l | NC | 20 |
| n-Butylbenzene | ND | ND | ug/l | NC | 20 |
| Dichlorodifluoromethane | ND | ND | ug/l | NC | 20 |
| Hexachlorobutadiene | ND | ND | ug/l | NC | 20 |
| Isopropylbenzene | ND | ND | ug/l | NC | 20 |
| p-Isopropyltoluene | ND | ND | ug/l | NC | 20 |
| Naphthalene | ND | ND | ug/l | NC | 20 |
| n-Propylbenzene | ND | ND | ug/l | NC | 20 |
| sec-Butylbenzene | ND | ND | ug/l | NC | 20 |
| tert-Butylbenzene | ND | ND | ug/l | NC | 20 |
| 1,2,3-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trichlorobenzene | ND | ND | ug/l | NC | 20 |
| 1,2,4-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| 1,3,5-Trimethylbenzene | ND | ND | ug/l | NC | 20 |
| Bromobenzene | ND | ND | ug/l | NC | 20 |
| o-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| p-Chlorotoluene | ND | ND | ug/l | NC | 20 |
| Dibromomethane | ND | ND | ug/l | NC | 20 |
| 1,2-Dibromoethane | ND | ND | ug/l | NC | 20 |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

| Parameter | Native Sample | Duplicate Sample | Units | RPD | RPD Limits |
|--|---------------------------------|-----------------------|-------------|-------------|----------------------|
| Volatile Organics by GC/MS - Westborough | Lab Associated sample(s): 01-09 | QC Batch ID: WG417697 | 7-4 QC Samp | le: L100881 | 2-01 Client ID: MW-3 |
| 1,2-Dibromo-3-chloropropane | ND | ND | ug/l | NC | 20 |
| 1,3-Dichloropropane | ND | ND | ug/l | NC | 20 |
| Methyl tert butyl ether | ND | ND | ug/l | NC | 20 |

| | | | | | Acceptance | |
|------------------------|-----------|-----------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | %Recovery | Qualifier | Criteria | |
| 1,2-Dichlorobenzene-d4 | 101 | | 101 | | 80-120 | |
| 4-Bromofluorobenzene | 96 | | 97 | | 80-120 | |

METALS



06/11/10

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

SAMPLE RESULTS

Lab ID: L1008812-01

Client ID: MW-3

Sample Location: WALPOLE, MA

Matrix: Water

Lab Number: L1008812

Report Date: 06/21/10

Date Collected: 06/11/10 08:02

Date Received:

Field Prep: See Narrative

| Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst | |
|--|--|------------------------|--|---|--|------------------|--|---|----------------------|---------|--|
| MCP Dissolved Metals - Westborough Lab | | | | | | | | | | | |
| ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 03:32 | EPA 3005A | 97,6020A | ВМ | |
| ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| 0.036 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | AI | |
| ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | AI | |
| ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | AI | |
| ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:34 | EPA 7470A | 97,7470A | EZ | |
| ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 14:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 03:32 | EPA 3005A | 97,6020A | ВМ | |
| ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | Al | |
| ND | | mg/l | 0.050 | | 1 | 06/14/10 13:10 | 06/16/10 17:57 | EPA 3005A | 97,6010B | AI | |
| | ND ND ND ND ND ND ND ND ND ND ND ND ND N | Als - Westborough Land | Als - Westborough Lab ND mg/l ND mg/l 0.036 mg/l ND mg/l | ND mg/l 0.0020 ND mg/l 0.005 0.036 mg/l 0.010 ND mg/l 0.004 ND mg/l 0.004 ND mg/l 0.004 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.0025 ND mg/l 0.0025 ND mg/l 0.010 ND mg/l 0.007 ND mg/l 0.0020 ND mg/l 0.0010 | ND mg/l 0.0020 ND mg/l 0.005 0.036 mg/l 0.010 ND mg/l 0.004 ND mg/l 0.004 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.0002 ND mg/l 0.0002 ND mg/l 0.010 ND mg/l 0.0002 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.010 ND mg/l 0.007 ND mg/l 0.0020 ND mg/l 0.0020 ND mg/l 0.0020 | ND | Result Qualifier Units RL MDL Factor Prepared Als - Westborough Lab ND mg/l 0.0020 4 06/12/10 10:30 ND mg/l 0.005 1 06/14/10 13:10 0.036 mg/l 0.010 1 06/14/10 13:10 ND mg/l 0.004 1 06/14/10 13:10 ND mg/l 0.01 1 06/14/10 13:10 ND mg/l 0.010 1 06/14/10 13:10 ND mg/l 0.0002 1 06/17/10 17:30 ND mg/l 0.0025 1 06/14/10 13:10 ND mg/l 0.007 1 06/14/10 13:10 ND mg/l 0.0020 4 06/12/10 10:30 ND mg/l 0.0020 4 06/12/10 10:30 ND mg/l 0.0020 < | Result Qualifier Units RL MDL Factor Prepared Analyzed Als - Westborough Lab ND mg/l 0.0020 4 06/12/10 10:30 06/15/10 03:32 ND mg/l 0.005 1 06/14/10 13:10 06/16/10 17:57 0.036 mg/l 0.010 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.004 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.004 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.01 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.010 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.002 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.025 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.010 1 06/14/10 13:10 06/16/10 17:57 ND mg/l 0.002 | ND | ND | |



L1008812

Project Name: WALPOLE PARK SOUTH Lab Number:

Project Number: 12700058-003 **Report Date:** 06/21/10

SAMPLE RESULTS

Lab ID: L1008812-02

Client ID: RIZ-3

Sample Location: WALPOLE, MA

Matrix: Water

Date Collected: 06/11/10 08:56

Date Received: 06/11/10

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-----------|-------------|-------|--------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Met | als - Wes | stborough L | ab | | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 0 06/15/10 03:38 | EPA 3005A | 97,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Barium, Dissolved | 0.161 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 0 06/18/10 11:35 | EPA 7470A | 97,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 0 06/17/10 15:00 | EPA 3005A | 97,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 0 06/15/10 03:38 | EPA 3005A | 97,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | | 1 | 06/14/10 13:10 | 0 06/16/10 18:00 | EPA 3005A | 97,6010B | Al |
| | | | | | | | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 Lab Number: **Report Date:**

L1008812 06/21/10

SAMPLE RESULTS

Lab ID:

L1008812-03

Client ID:

MW-2

Sample Location:

WALPOLE, MA

Matrix:

Water

Date Collected:

06/11/10 09:33

Date Received: Field Prep:

06/11/10

See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-----------|------------|-------|--------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Meta | als - Wes | tborough L | ab | | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 03:44 | EPA 3005A | 97,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Barium, Dissolved | 0.070 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:37 | EPA 7470A | 97,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 15:04 | EPA 3005A | 97,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 03:44 | EPA 3005A | 97,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | | 1 | 06/14/10 13:10 | 06/16/10 18:04 | EPA 3005A | 97,6010B | Al |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

Lab ID: L1008812-04

Client ID: GHC-6

Date Collected:
Date Received:

06/11/10 10:08

Sample Location:

WALPOLE, MA

06/11/10

Matrix:

Water

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-----------|------------|-------|--------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Met | als - Wes | tborough L | ab | | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:09 | EPA 3005A | 97,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | AI |
| Barium, Dissolved | 0.063 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Lead, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:42 | EPA 7470A | 97,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Selenium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 14:34 | EPA 3005A | 97,6010B | Al |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:09 | EPA 3005A | 97,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:24 | EPA 3005A | 97,6010B | Al |
| Zinc, Dissolved | ND | | mg/l | 0.050 | | 1 | | 06/16/10 17:24 | | 97,6010B | Al |
| | | | | | | | | | | | |

SAMPLE RESULTS



Project Name: WALPOLE PARK SOUTH

12700058-003

Lab Number: **Report Date:**

L1008812

06/21/10

Lab ID: L1008812-05

Client ID: RIZ-9

Sample Location: WALPOLE, MA

Matrix: Water

Project Number:

Date Collected:

06/11/10 11:07

Date Received:

06/14/10 13:10 06/16/10 17:34 EPA 3005A

06/11/10

Field Prep:

See Narrative

Dilution Date Date Prep Analytical Method Method Factor **Prepared Analyzed** Qualifier Units RL MDL **Parameter** Result **Analyst** MCP Dissolved Metals - Westborough Lab Antimony, Dissolved ND mg/l 0.0020 4 06/12/10 10:30 06/15/10 04:27 EPA 3005A 97,6020A ВМ ND 1 97,6010B Arsenic, Dissolved mg/l 0.005 06/14/10 13:10 06/16/10 17:34 EPA 3005A ΑI ND 1 0.010 97,6010B Barium, Dissolved mg/l 06/14/10 13:10 06/16/10 17:34 EPA 3005A ΑI Beryllium, Dissolved ND mg/l 0.004 1 06/14/10 13:10 06/16/10 17:34 EPA 3005A 97,6010B ΑI ND 0.004 1 97,6010B Cadmium, Dissolved 06/14/10 13:10 06/16/10 17:34 EPA 3005A ΑI mg/l ND 0.01 1 97,6010B Chromium, Dissolved mg/l 06/14/10 13:10 06/16/10 17:34 EPA 3005A ΑI Lead, Dissolved ND 0.010 --1 06/14/10 13:10 06/16/10 17:34 EPA 3005A 97,6010B ΑI mg/l Mercury, Dissolved ND mg/l 0.0002 1 06/17/10 17:30 06/18/10 11:46 EPA 7470A 97,7470A ΕZ ND 1 97,6010B Nickel, Dissolved 0.025 06/14/10 13:10 06/16/10 17:34 EPA 3005A ΑI mg/l Selenium, Dissolved ND 0.010 --1 06/14/10 13:10 06/16/10 17:34 EPA 3005A 97,6010B ΑI mg/l Silver, Dissolved ND 0.007 1 06/14/10 13:10 06/17/10 14:44 EPA 3005A 97,6010B mg/l ΑI --Thallium, Dissolved ND 0.0020 4 06/12/10 10:30 06/15/10 04:27 EPA 3005A 97,6020A BM mg/l --Vanadium, Dissolved ND 0.010 1 06/14/10 13:10 06/16/10 17:34 EPA 3005A 97,6010B mg/l ΑI ND 0.050 1 97.6010B ΑI

mg/l

SAMPLE RESULTS



Zinc, Dissolved

Project Name: WALPOLE PARK SOUTH Lab Number:

Report Date:

L1008812

06/21/10

Project Number: 1270

12700058-003

SAMPLE RESULTS

06/11/10 12:30

Lab ID: Client ID: L1008812-06

RIZ-10

Date Collected: 06/11/10
Date Received: 06/11/10

Sample Location:

WALPOLE, MA

Field Prep: See Narrative

Matrix:

Water

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-----------|-------------|-------|--------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Meta | als - Wes | tborough La | ab | | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:33 | EPA 3005A | 97,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Barium, Dissolved | 0.107 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Chromium, Dissolved | ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:48 | EPA 7470A | 97,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Silver, Dissolved | ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 14:47 | EPA 3005A | 97,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:33 | EPA 3005A | 97,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:37 | EPA 3005A | 97,6010B | AI |
| Zinc, Dissolved | ND | | mg/l | 0.050 | | 1 | | 06/16/10 17:37 | | 97,6010B | Al |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: Report Date: L1008812

06/21/10

SAMPLE RESULTS

mg/l

0.005

0.010

0.004

0.004

0.01

0.010

0.0002

0.025

0.010

0.007

0.0020

0.010

0.050

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L1008812-07

Client ID: RIZ-8

Sample Location: WALPOLE, MA

ND

Matrix: Water

Lab ID:

Arsenic, Dissolved

Barium, Dissolved

Beryllium, Dissolved

Cadmium, Dissolved

Chromium, Dissolved

Lead, Dissolved

Mercury, Dissolved

Selenium, Dissolved

Thallium, Dissolved

Vanadium, Dissolved

Nickel, Dissolved

Silver, Dissolved

Zinc, Dissolved

Date Collected:

06/11/10 13:40

Date Received:

06/14/10 13:10 06/16/10 17:50 EPA 3005A

06/17/10 17:30 06/18/10 11:49 EPA 7470A

06/14/10 13:10 06/16/10 17:50 EPA 3005A

06/14/10 13:10 06/16/10 17:50 EPA 3005A

06/14/10 13:10 06/17/10 14:51 EPA 3005A

06/12/10 10:30 06/15/10 04:39 EPA 3005A

06/14/10 13:10 06/16/10 17:50 EPA 3005A

06/14/10 13:10 06/16/10 17:50 EPA 3005A

06/11/10

Field Prep:

See Narrative

97,6010B

97,6010B

97,6010B

97,6010B

97,6010B

97,6010B

97,7470A

97,6010B

97,6010B

97,6010B

97,6020A

97,6010B

97.6010B

ΑI

ΑI

ΑI

ΑI

ΑI

ΑI

ΕZ

ΑI

ΑI

ΑI

BM

ΑI

ΑI

Dilution Date Date Prep Analytical Method **Prepared** Method Factor **Analyzed** Qualifier Units RL MDL **Parameter** Result **Analyst** MCP Dissolved Metals - Westborough Lab Antimony, Dissolved ND mg/l 0.0020 4 06/12/10 10:30 06/15/10 04:39 EPA 3005A 97,6020A BM

1

1

1

1

1

1

1

1

1

1

4

1

1

ALPHA

Project Name: Lab Number: WALPOLE PARK SOUTH 06/21/10

Project Number: 12700058-003

Report Date:

L1008812

SAMPLE RESULTS

Lab ID: L1008812-08

Client ID: MW-9

Sample Location: WALPOLE, MA

Matrix: Water Date Collected: 06/11/10 14:10

Date Received: 06/11/10

Field Prep: See Narrative

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Prep Method | Analytical Method | Analyst |
|----------------------|-----------|-------------|-------|--------|-----|--------------------|------------------|------------------|----------------|----------------------|---------|
| MCP Dissolved Met | als - Wes | stborough L | .ab | | | | | | | | |
| Antimony, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:45 | EPA 3005A | 97,6020A | ВМ |
| Arsenic, Dissolved | ND | | mg/l | 0.005 | | 1 | 06/14/10 13:10 |) 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Barium, Dissolved | 0.064 | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Beryllium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Cadmium, Dissolved | ND | | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | Al |
| Chromium, Dissolved | ND | | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Lead, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Mercury, Dissolved | ND | | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:51 | EPA 7470A | 97,7470A | EZ |
| Nickel, Dissolved | ND | | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Selenium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | Al |
| Silver, Dissolved | ND | | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 14:54 | EPA 3005A | 97,6010B | AI |
| Thallium, Dissolved | ND | | mg/l | 0.0020 | | 4 | 06/12/10 10:30 | 06/15/10 04:45 | EPA 3005A | 97,6020A | ВМ |
| Vanadium, Dissolved | ND | | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| Zinc, Dissolved | 0.051 | | mg/l | 0.050 | | 1 | 06/14/10 13:10 |) 06/16/10 17:54 | EPA 3005A | 97,6010B | AI |
| | | | | | | | | | | | |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date: 06/21/10

Method Blank Analysis Batch Quality Control

| Parameter | Result Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|------------------------|---------------------|----------|---------|-------|--------------------|------------------|------------------|----------------------|----|
| MCP Dissolved Metals - | Westborough Lab for | sample(s |): 01-0 | 8 Bat | ch: WG417 | '601-1 | | | |
| Antimony, Dissolved | ND | mg/l | 0.0005 | | 1 | 06/12/10 10:30 | 06/14/10 18:34 | 97,6020A | ВМ |
| Thallium, Dissolved | ND | mg/l | 0.0005 | | 1 | 06/12/10 10:30 | 06/14/10 18:34 | 97,6020A | BM |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | Analyst |
|----------------------|-----------------------|-----------|--------|------|--------------------|------------------|------------------|----------------------|---------|
| MCP Dissolved Metals | - Westborough Lab for | sample(s) | : 01-0 | Bate | ch: WG417 | 7803-1 | | | |
| Arsenic, Dissolved | ND | mg/l | 0.005 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Barium, Dissolved | ND | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Beryllium, Dissolved | ND | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Cadmium, Dissolved | ND | mg/l | 0.004 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Chromium, Dissolved | ND | mg/l | 0.01 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Lead, Dissolved | ND | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Nickel, Dissolved | ND | mg/l | 0.025 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Selenium, Dissolved | ND | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Silver, Dissolved | ND | mg/l | 0.007 | | 1 | 06/14/10 13:10 | 06/17/10 14:07 | 97,6010B | Al |
| Vanadium, Dissolved | ND | mg/l | 0.010 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |
| Zinc, Dissolved | ND | mg/l | 0.050 | | 1 | 06/14/10 13:10 | 06/16/10 17:15 | 97,6010B | Al |

Prep Information

Digestion Method: EPA 3005A

| Parameter | Result Qualifier | Units | RL | MDL | Dilution Factor | Date Prepared | Date Analyzed | Analytical Method | |
|---------------------|-------------------------|----------|-----------|-----|--------------------|------------------|------------------|----------------------|----|
| MCP Dissolved Metal | s - Westborough Lab for | sample(s | s): 01-08 | Bat | ch: WG418 | 3542-1 | | | |
| Mercury, Dissolved | ND | mg/l | 0.0002 | | 1 | 06/17/10 17:30 | 06/18/10 11:28 | 97,7470A | EZ |



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 **Report Date:** 06/21/10

Method Blank Analysis Batch Quality Control

Prep Information

Digestion Method: EPA 7470A



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number:

L1008812

Report Date:

06/21/10

| Parameter | LCS %Recovery (| Qual % | LCSD Recovery | / Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|--------------------|-----------|------------------|------------|---------------------|-----|------|------------|
| MCP Dissolved Metals - Westborough Lab | Associated sample(| s): 01-08 | Batch: | WG417601-2 | WG417601-3 | | | |
| Antimony, Dissolved | 95 | | 96 | | 80-120 | 1 | | 20 |
| Thallium, Dissolved | 96 | | 96 | | 80-120 | 0 | | 20 |
| MCP Dissolved Metals - Westborough Lab | Associated sample(| s): 01-08 | Batch: | WG417803-2 | WG417803-3 | | | |
| Arsenic, Dissolved | 114 | | 115 | | 80-120 | 1 | | 20 |
| Barium, Dissolved | 98 | | 100 | | 80-120 | 2 | | 20 |
| Beryllium, Dissolved | 100 | | 102 | | 80-120 | 2 | | 20 |
| Cadmium, Dissolved | 109 | | 110 | | 80-120 | 1 | | 20 |
| Chromium, Dissolved | 95 | | 100 | | 80-120 | 5 | | 20 |
| Lead, Dissolved | 107 | | 110 | | 80-120 | 3 | | 20 |
| Nickel, Dissolved | 98 | | 99 | | 80-120 | 1 | | 20 |
| Selenium, Dissolved | 113 | | 115 | | 80-120 | 2 | | 20 |
| Silver, Dissolved | 92 | | 94 | | 80-120 | 2 | | 20 |
| Vanadium, Dissolved | 101 | | 102 | | 80-120 | 1 | | 20 |
| Zinc, Dissolved | 101 | | 102 | | 80-120 | 1 | | 20 |
| MCP Dissolved Metals - Westborough Lab | Associated sample(| s): 01-08 | Batch: | WG418542-2 | WG418542-3 | | | |
| Mercury, Dissolved | 114 | | 111 | | 80-120 | 3 | | 20 |



Matrix Spike Analysis Batch Quality Control

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812

Report Date: 06/21/10

| Parameter | Native Sample | MS Added | MS Found %F | MS Recovery | MSD Qual Found | MSD %Recovery Qu | Recovery | / RPD Qı | RPD _{ual} Limits |
|--------------------------|-------------------|--------------|-----------------|----------------|-------------------|---------------------|-------------|-------------|------------------------------|
| MCP Dissolved Metals - V | Vestborough Lab A | Associated s | ample(s): 01-08 | QC Bat | ch ID: WG417601-4 | QC Sample: L | .1008812-04 | Client ID: | GHC-6 |
| Antimony, Dissolved | ND | 0.5 | 0.4899 | 98 | | - | 75-125 | - | 20 |
| Thallium, Dissolved | ND | 0.12 | 0.1159 | 96 | - | - | 75-125 | - | 20 |
| MCP Dissolved Metals - V | Vestborough Lab / | Associated s | ample(s): 01-08 | 3 QC Bat | ch ID: WG417803-4 | QC Sample: L | .1008812-04 | Client ID: | GHC-6 |
| Arsenic, Dissolved | ND | 0.12 | 0.140 | 117 | - | - | 75-125 | - | 20 |
| Barium, Dissolved | 0.063 | 2 | 2.04 | 99 | - | - | 75-125 | - | 20 |
| Beryllium, Dissolved | ND | 0.05 | 0.050 | 100 | - | - | 75-125 | - | 20 |
| Cadmium, Dissolved | ND | 0.051 | 0.055 | 108 | - | - | 75-125 | - | 20 |
| Chromium, Dissolved | ND | 0.2 | 0.20 | 100 | - | - | 75-125 | - | 20 |
| Lead, Dissolved | ND | 0.51 | 0.533 | 104 | - | - | 75-125 | - | 20 |
| Nickel, Dissolved | ND | 0.5 | 0.472 | 94 | - | - | 75-125 | - | 20 |
| Selenium, Dissolved | ND | 0.12 | 0.137 | 114 | - | - | 75-125 | - | 20 |
| Silver, Dissolved | ND | 0.05 | 0.047 | 95 | - | - | 75-125 | - | 20 |
| Vanadium, Dissolved | ND | 0.5 | 0.531 | 106 | - | - | 75-125 | - | 20 |
| Zinc, Dissolved | ND | 0.5 | 0.511 | 102 | - | - | 75-125 | - | 20 |
| MCP Dissolved Metals - V | Vestborough Lab / | Associated s | ample(s): 01-08 | 3 QC Bat | ch ID: WG418542-4 | QC Sample: L | .1008812-04 | Client ID: | GHC-6 |
| Mercury, Dissolved | ND | 0.001 | 0.0011 | 115 | - | | 75-125 | - | 20 |

Lab Number: L1008812

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 **Report Date:** 06/21/10

Sample Receipt and Container Information

Were project specific reporting limits specified?

Reagent H2O Preserved Vials Frozen on: NA

Cooler Information Custody Seal

Cooler

A Absent

| Container Info | ormation | | | Temp | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|---|
| Container ID | Container Type | Cooler | рΗ | deg C | Pres | Seal | Analysis(*) |
| L1008812-01A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-01B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-01C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-TL-6020S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-PB-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-02A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-02B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-02C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-AS-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BB-6010S-10(180),MCP-PB-6010S-10(180),MCP-PB-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-03A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-03B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |



Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003

Lab Number: L1008812 **Report Date:** 06/21/10

| Container Info | ormation | | | Temp | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|--|
| Container ID | Container Type | Cooler | рН | deg C | Pres | Seal | Analysis(*) |
| L1008812-03C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-04A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-04B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-04C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-PB-6010S-10(180),MCP-PB-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-04D | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-PB-6010S-10(180),MCP-PB-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-05A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-05B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-05C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-AS-6010S-10(180),MCP-AS-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-PB-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-06A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |



Lab Number: L1008812

Project Name: WALPOLE PARK SOUTH

Project Number: 12700058-003 **Report Date:** 06/21/10

| Container Info | rmation | | | Temp | | | |
|----------------|----------------------------------|--------|-----|-------|------|--------|---|
| Container ID | Container Type | Cooler | рН | deg C | Pres | Seal | Analysis(*) |
| L1008812-06B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-06C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-XN-6010S-10(180),MCP-AS-6010S-10(180),MCP-TL-6020S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-07A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-07B | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-07C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-08A | Vial Ascorbic Acid/HCl preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-08B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-08C | Plastic 250ml HNO3 preserved | A | <2 | 4 | Y | Absent | MCP-CD-6010S-10(180),MCP-7470S-10(28),MCP-AG-6010S-10(180),MCP-SB-6020S-10(180),MCP-ZN-6010S-10(180),MCP-AS-6010S-10(180),MCP-CR-6010S-10(180),MCP-BA-6010S-10(180),MCP-BE-6010S-10(180),MCP-BE-6010S-10(180),MCP-BB-6010S-10(180),MCP-NI-6010S-10(180),MCP-NI-6010S-10(180),MCP-SE-6010S-10(180),MCP-SE-6010S-10(180),MCP-V-6010S-10(180) |
| L1008812-09A | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | 524.2(14) |
| L1008812-09B | Vial Ascorbic Acid/HCI preserved | Α | N/A | 4 | Υ | Absent | - |

Container Comments

L1008812-01C

L1008812-07C



Project Name:WALPOLE PARK SOUTHLab Number:L1008812Project Number:12700058-003Report Date:06/21/10

GLOSSARY

Acronyms

EPA · Environmental Protection Agency.

 LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD · Laboratory Control Sample Duplicate: Refer to LCS.

MDL • Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

MS • Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available.

MSD · Matrix Spike Sample Duplicate: Refer to MS.

NA · Not Applicable.

NC • Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

NI · Not Ignitable.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Data Qualifiers

- A Spectra identified as "Aldol Condensation Product".
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than five times (5x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- ${\bf E} \qquad \hbox{-Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.}$
- The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- The quality control sample exceeds the associated acceptance criteria. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.

Report Format: Data Usability Report



Project Name:WALPOLE PARK SOUTHLab Number:L1008812Project Number:12700058-003Report Date:06/21/10

Data Qualifiers

J Estimated value. This represents an estimated concentration for Tentatively Identified Compounds (TICs).

 ${\bf ND}$ • Not detected at the reporting limit (RL) for the sample.

Report Format: Data Usability Report



Project Name: WALPOLE PARK SOUTH Lab Number: L1008812

Project Number: 12700058-003 Report Date: 06/21/10

REFERENCES

Methods for the Determination of Organic Compounds in Drinking Water - Supplement II. EPA/600/R-92/129, August 1992.

97 EPA Test Methods (SW-846) with QC Requirements & Performance Standards for the Analysis of EPA SW-846 Methods under the Massachusetts Contingency Plan, WSC-CAM-IIA, IIB, IIIA, IIIB, IIIC, IIID, VA, VB, VC, VIA, VIB, VIIIA and VIIIB, July 2010.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certificate/Approval Program Summary

Last revised June 17, 2010 - Westboro Facility

The following list includes only those analytes/methods for which certification/approval is currently held. For a complete listing of analytes for the referenced methods, please contact your Alpha Customer Service Representative.

Connecticut Department of Public Health Certificate/Lab ID: PH-0574. NELAP Accredited Solid Waste/Soil.

Drinking Water (Inorganic Parameters: Color, pH, Turbidity, Conductivity, Alkalinity, Chloride, Free Residual Chlorine, Fluoride, Calcium Hardness, Sulfate, Nitrate, Nitrite, Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Vanadium, Zinc, Total Dissolved Solids, Total Organic Carbon, Total Cyanide, Perchlorate.

Organic Parameters: Haloacetic Acids, Volatile Organics 524.2, Total Trihalomethanes 524.2, 1,2-Dibromo-3-chloropropane (DBCP), Ethylene Dibromide (EDB).)

Wastewater/Non-Potable Water (Inorganic Parameters: Color, pH, Conductivity, Acidity, Alkalinity, Chloride, Total Residual Chlorine, Fluoride, Total Hardness, Calcium Hardness, Silica, Sulfate, Sulfide, Ammonia, Kjeldahl Nitrogen, Nitrate, Nitrite, O-Phosphate, Total Phosphorus, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Strontium, Thallium, Tin, Titanium, Vanadium, Zinc, Total Residue (Solids), Total Dissolved Solids, Total Suspended Solids (non-filterable), BOD, CBOD, COD, TOC, Total Cyanide, Phenolics, Foaming Agents (MBAS), Bromide, Oil and Grease. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Acid Extractables (Phenols), Benzidines, Phthalate Esters, Nitrosamines, Nitroaromatics & Isophorone, Polynuclear Aromatic Hydrocarbons, Haloethers, Chlorinated Hydrocarbons, Volatile Organics, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH.) Solid Waste/Soil (Inorganic Parameters: Lead in Paint, pH, Aluminum, Antimony, Arsenic, Barium, Beryllium, Boron, Cadmium, Calcium, Chromium, Hexavalent Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Mercury, Molybdenum, Nickel, Potassium, Selenium, Silver, Sodium, Thallium, Tin, Vanadium, Zinc, Total Cyanide, Ignitability, Phenolics, Corrosivity, TCLP Leach (1311), Reactivity. Organic Parameters: PCBs, Organochlorine Pesticides, Technical Chlordane, Toxaphene, Extractable Petroleum Hydrocarbons (ETPH), MA-EPH, MA-VPH, Dicamba, 2,4-D, 2,4,5-T, 2,4,5-TP(Silvex), Volatile Organics, Acid Extractables (Phenols), 3.3'-Dichlorobenzidine, Phthalates,

Maine Department of Human Services Certificate/Lab ID: 2009024.

Drinking Water (Inorganic Parameters: SM9215B, 9221E, 9222B, 9222D, 9223B, EPA 180.1, 300.0, 353.2, SM2130B, 2320B, 4500Cl-D, 4500CN-C, 4500CN-E, 4500F-C, 4500H+B,4500NO3-F, EPA 200.7, EPA 200.8, 245.1. Organic Parameters: 504.1, 524.2, SM 6251B.)

Wastewater/Non-Potable Water (Inorganic Parameters: EPA 120.1, 1664A, 350.1, 351.1, 353.2, 410.4, 420.1, Lachat 10-107-06-1-B, SM2320B, 2340B, 2510B, 2540C, 2540D, 426C, 4500Cl-D, 4500Cl-E, 4500CN-C, 4500CN-E, 4500F-B, 4500F-C, 4500H+B, 4500Norg-B, 4500Norg-C, 4500NH3-B, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500P-B.5, 4500P-E, 5210B, 5220D, 5310C, EPA 200.7, 200.8, 245.1. Organic Parameters: 608, 624.)

Massachusetts Department of Environmental Protection Certificate/Lab ID: M-MA086.

Nitrosamines, Nitroaromatics & Cyclic Ketones, PAHs, Haloethers, Chlorinated Hydrocarbons.)

Drinking Water

Inorganic Parameters: (EPA 200.8 for: Sb,As,Ba,Be,Cd,Cr,Cu,Pb,Ni,Se,Tl)

(EPA 200.7 for: Ba,Be,Ca,Cd,Cr,Cu,Na,Ni) 245.1, (300.0 for: Nitrate-N, Fluoride, Sulfate)

353.2 for: Nitrate-N, Nitrite-N; SM4500NO3-F, 4500F-C, 4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, 2320B, SM2540C. SM4500H-B.

Organic Parameters: (EPA 524.2 for: Trihalomethanes, Volatile Organics)

(504.1 for: 1,2-Dibromoethane, 1,2-Dibromo-3-Chloropropane), 314.0, 332.

Microbiology Parameters: SM9215B; ENZ. SUB. SM9223; MF-SM9222D

Non-Potable Water

Inorganic Parameters:, (EPA 200.8 for: Al,Sb,As,Be,Cd,Cr,Cu,Pb,Mn,Ni,Se,Ag,Tl,Zn)

(EPA 200.7 for: Al,Sb,As,Be,Cd,Cr,Co,Cu,Fe,Pb,Mn,Mo,Ni,Se,Aq,Sr,Ti,Tl, V,Zn,Ca,Mq,Na,K)

245.1, SM4500H,B, EPA 120.1, SM2510B, 2540C, 2540B, 2340B, 2320B, 4500CL-E, 4500F-BC, 426C, SM4500NH3-BH, (EPA 350.1 for: Ammonia-N), LACHAT 10-107-06-1-B for Ammonia-N, SM4500NO3-F, 353.2 for Nitrate-N, SM4500NH3-B,C-Titr, SM4500NH3-BC-NES, EPA 351.1, SM4500P-E, 4500P-B,E, 5220D, EPA 410.4, SM 5210B, 5310C, 4500CL-D, EPA 1664, SM14 510AC, EPA 420, SM4500-CN-CE, SM2540D.

Organic Parameters: (EPA 624 for Volatile Halocarbons, Volatile Aromatics)

(608 for: Chlordane, Aldrin, Dieldrin, DDD, DDE, DDT, Heptachlor, Heptachlor Epoxide, PCBs-Water), EPA 625 for SVOC Acid Extractables and SVOC Base/Neutral Extractables, 600/4-81-045-PCB-Oil

New Hampshire Department of Environmental Services Certificate/Lab ID: 200307. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM6215B, 9222B, 9223B Colilert, EPA 200.7, 200.8, 245.2, 120.1, 300.0, 314.0, SM4500CN-E, 4500H+B, 4500NO3-F, 2320B, 2510B, 2540C, 4500F-C, 5310C, 2120B, EPA 331.0. Organic Parameters: 504.1, 524.2, SM6251B.)

Non-Potable Water (Inorganic Parameters: SM9222D, 9221B, 9222B, 9221E-EC, EPA 200.7, 200.8, 245.1, 245.2, SW-846 6010B, 6020, 7196A, 7470A, SM3500-CR-D, EPA 120.1, 300.0, 350.1, 351.1, 353.2, 420.1, 1664A, SW-846 9010, 9030, 9040B, SM426C, SM2310B, 2540B, 2540D, 4500H+B, 4500NH3-H, 4500NH3-E, 4500NO2-B, 4500P-E, 4500-S2-D, 5210B, 2320B, 2540C, 4500F-C, 5310C, 5540C, LACHAT 10-117-07-1-B, LACHAT 10-107-06-1-B, LACHAT 10-107-04-1-J, LACHAT 10-117-07-1-A, SM4500CL-E, LACHAT 10-204-00-1-A, LACHAT 10-107-06-2-D. Organic Parameters: SW-846 3005A, 3015A, 3510C, 5030B, 8021B, 8260B, 8270C, 8330, EPA 624, 625, 608, SW-846 8082, 8081A.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 6010B, 7196A, 7471A, 7.3.3.2, 7.3.4.2, 1010, 1030, 9010, 9012A, 9014, 9030B, 9040, 9045C, 9050C, 1311, 3005A, 3050B, 3051A. Organic Parameters: SW-846 3540C, 3545, 3580A, 5030B, 5035, 8021B, 8260B, 8270C, 8330, 8151A, 8082, 8081A.)

New Jersey Department of Environmental Protection Certificate/Lab ID: MA935. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9222B, 9221E, 9223B, 9215B, 4500NO3-F, 4500F-C, EPA 300.0, 200.7, 2540C, 2320B, 314.0, SM2120B, 2510B, 5310C, SM4500H-B, EPA 200.8, 245.2. Organic Parameters: 504.1, SM6251B, 524.2.)

Non-Potable Water (Inorganic Parameters: SM5210B, EPA 410.4, SM5220D, 4500Cl-D, EPA 300.0, SM2120B, SM4500F-BC, EPA 200.7, 351.1, LACHAT 10-107-06-2-D, EPA 353.2, SM4500NO3-F, 4500NO2-B, EPA 1664A, SM5310B, C or D, 4500-PE, EPA 420.1, SM4500P-B5+E, 2540B, 2540C, 2540D, EPA 120.1, SM2510B, SM15 426C, SM9221CE, 9222D, 9221B, 9222B, 9215B, 2310B, 2320B, 4500NH3-H, 4500-S D, EPA 350.1, SM5210B, SW-846 3015, 6020, 7470A, 5540C, 4500H-B, EPA 200.8, SM3500Cr-D, EPA 245.1, 245.2, SW-846 9040B, 3005A, EPA 6010B, 7196A, SW-846 9010B, 9030B. Organic Parameters: SW-846 8260B, 8270C, 3510C, EPA 608, 624, 625, SW-846 5030B, 8021B, 8081A, 8082, 8151A, 8330, NJ OQA-QAM-025 Rev.7.)

Solid & Chemical Materials (Inorganic Parameters: SW-846 9040B, 3005A, 6010B, 7196A, 5030B, 9010B, 9030B, 1030, 1311, 3050B, 3051, 7471A, 9014, 9012A, 9045C, 9050A, 9065. Organic Parameters: SW-846 8021B, 8081A, 8082, 8151A, 8330, 8260B, 8270C, 1311, 1312, 3540C, 3545, 3550B, 3580A, 5035L, 5035H, NJ OQA-QAM-025 Rev.7.)

New York Department of Health Certificate/Lab ID: 11148. NELAP Accredited.

Drinking Water (Inorganic Parameters: SM9223B, 9222B, 9215B, EPA 200.8, 200.7, 245.2, SM5310C, EPA 314.0, 332.0, SM2320B, EPA 300.0, SM2120B, 4500CN-E, 4500F-C, 4500H-B, 4500NO3-F, 2540C, EPA 120.1, SM 2510B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: SM9221E, 9222D, 9221B, 9222B, 9215B, 5210B, EPA 410.4, SM5220D, 2310B-4a, 2320B, EPA 200.7, 300.0, LACHAT 10-117-07-1A or B, SM4500Cl-E, 4500F-C, SM15 426C, EPA 350.1, LACHAT 10-107-06-1-B, SM4500NH3-H, EPA 351.1, LACHAT 10-107-06-2, EPA 353.2, LACHAT 10-107-041-C, SM4500-NO3-F, 4500-NO2-B, 4500P-E, 2540C, 2540B, 2540D, EPA 200.8, EPA 6010B, 6020, EPA 7196A, S\M3500Cr-D, EPA 245.1, 245.2, 7470A, SM2120B, SM4500-CN-E LACHAT 10-204-00-1-A, EPA 9040B, SM4500-HB, EPA 1664A, SM5310C, EPA 420.1, SM14 510C, EPA 120.1, SM2510B, SM4500S-D, SM5540C, EPA 3005A, 3015. Organic Parameters: EPA 624, 8260B, 8270C, 625, 608, 8081A, 8151A, 8330, 8082, EPA 3510C, 5030B, 9010B, 9030B.)

Solid & Hazardous Waste (Inorganic Parameters: 1010, 1030, SW-846 Ch 7 Sec 7.3, EPA 6010B, 7196A, 7471A, 9012A, 9014, 9040B, 9045C, 9065, 9050, EPA 1311, 1312, 3005A, 3050B, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8081A, 8151A, 8330, 8082, 3540C, 3545, 3546, 3580, 5030B, 5035.)

North Carolina Department of the Environment and Natural Resources <u>Certificate/Lab ID</u>: 666. <u>Organic Parameters</u>: MA-EPH, MA-VPH.

Pennsylvania Department of Environmental Protection Certificate/Lab ID: 68-03671. *NELAP Accredited. Non-Potable Water* (Organic Parameters: EPA 3510C, 5030B, 625, 624. 608, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1010, 1030, 1311, 3050B, 3051, 6010B, EPA 7.3.3.2, EPA 7.3.4.2, 7196A, 7471A, 9010B, 9012A, 9014, 9040B, 9045C, 9050, 9065. Organic Parameters: 3540C, 3545, 3580A, 5035, 8021B, 8081A, 8082, 8151A, 8260B, 8270C, 8330)

Rhode Island Department of Health Certificate/Lab ID: LAO00065. *NELAP Accredited via NY-DOH.*Refer to MA-DEP Certificate for Potable and Non-Potable Water.
Refer to NY-DOH Certificate for Potable and Non-Potable Water.

Texas Commisson on Environmental Quality <u>Certificate/Lab ID</u>: T104704476-09-1. **NELAP Accredited.** Non-Potable Water (<u>Inorganic Parameters</u>: EPA 120.1, 1664, 200.7, 200.8, 245.1, 245.2, 300.0, 350.1, 351.1, 353.2, 376.2, 410.4, 420.1, 6010, 6020, 7196, 7470, 9040, SM 2120B, 2310B, 2320B, 2510B, 2540B, 2540C, 2540D, 426C, 4500CL-E, 4500CN-E, 4500F-C, 4500H+B, 4500NH3-H, 4500NO2B, 4500P-E, 4500 S2⁻ D, 510C, 5210B, 5220D, 5310C, 5540C. Organic Parameters: EPA 608, 624, 625, 8081, 8082, 8151, 8260, 8270, 8330.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 1311, 1312, 9012, 9014, 9040, 9045, 9050, 9065.)

Department of Defense Certificate/Lab ID: L2217.

Drinking Water (Inorganic Parameters: SM 4500H-B. Organic Parameters: EPA 524.2, 504.1.)

Non-Potable Water (Inorganic Parameters: EPA 200.7, 200.8, 6010B, 6020, 245.1, 245.2, 7470A, 9040B, 300.0, 9251, 9038, 350.1, 353.2, 351.1, 314, 120.1, 9050A, 410.4, 9060, 1664, 420.1, LACHAT 10-107-06-1-B, SM 4500CN-E, 4500H-B, 4500CL-E, 4500F-BC, 4500SO4-E, 426C, 4500NH3-B, 4500NH3-H, 4500NO3-F, 4500NO2-B, 4500Norg-C, 4500PE, 2510B, 5540C, 5220D, 5310C, 2540B, 2540C, 2540D, 510C, 4500S2-AD, 3005A, 3015, 9010B, 9030B. Organic Parameters: EPA 8260B, 8270C, 8330, 625, 8082, 8151A, 8081A, 3510C, 5030B, MassDEP EPH, MassDEP VPH.)

Solid & Hazardous Waste (Inorganic Parameters: EPA 200.7, 6010B, 7471A, 9040B, 9045C, 9065, 420.1, 9012A, 6860, 1311, 1312, 3050B, 9030B, 3051, 9010B, 3540C, SM 510ABC, 4500CN-CE, 2540G, SW-846 7.3, Organic Parameters: EPA 8260B, 8270C, 8330, 8082, 8081A, 8151A, 3545, 3546, 3580, 5035, MassDEP EPH, MassDEP VPH.)

Analytes Not Accredited by NELAP

Certification is not available by NELAP for the following analytes: **EPA 8260B:** Freon-113, 1,2,4,5-Tetramethylbenzene, 4-Ethyltoluene. **EPA 8330A:** PETN, Picric Acid, Nitroglycerine, 2,6-DANT, 2,4-DANT. **EPA 8270C:** Methyl naphthalene, Dimethyl naphthalene, Total Methylnapthalenes, Total Dimethylnaphthalenes, 1,4-Diphenylhydrazine (Azobenzene). **EPA 625:** 4-Chloroaniline. **EPA 350.1** for Ammonia in a Soil matrix.

| FORM NO: 01-01 (rev. 14-OCT-07) | MA MCP or CT RCP? | PLEASE ANSWER QUESTIONS ABOVE: | 9 201006/1- Trip Black | 8 MW-9 | 7 812-8 | 6 R1Z-10 | S RIZ-9 | F GHC-6 | 2 mw-2 | Q R12-3 | 8812.1 MW-3 | (Lab Use Only) Sample in | | | Other Project Specific Requirements/Comments/Detection Limits | These samples have been previously analyzed by Alpha | Email: Jan. Cannan Whetertech. 100 | Fax: 508 903 -2001 | EPhone: 56 9 903- 20 39 | Francisher int | Address One Grant St | Client Tetra Tech Rizzo | Client Information | TEL: 508-898-9220 TEL: 508-822-3300 FAX: 508-898-9193 FAX: 508-822-3288 | | CHAIN OF |
|---------------------------------|----------------------------------|--|------------------------|----------|---------|----------|---------|-------------------|--------|--|--------------|--------------------------|------------------------|--|---|--|------------------------------------|--|--------------------------------------|--|--|-----------------------------|------------------------------|---|--|--------------------------------|
| | Remoderied By: | | 6/9/10 | 0111 | 0481 | | 167 | 8001 | 0933 | 1 085% | 6/11/10 0802 | Date Time | ollectio | | ments/Detection Limits: | (0)(8)(0) | - | X Standard T RUSH (new confirmed fine) | Turn-Around Time | ALPHA Quote #: | Project Manager hay Junson for Camer | Project # 12700058 003 | Project Location: Walpole MA | Project Name: Whole Pour South | Project Information | CUSTODY |
| 1 | Date/Time | Container Type V P Preservative A(U) N | 15 Agric X | V | | | | | | | 6w KX | Initials | Sample Sampler's | | ₽Va 7 | | ellectory | appropriate | × | j. | | 234 | | | Re | PAGE OF Dai |
| | Received By: | 7 | | | | | | | | | × | / d/ / /see / /see | 1 | W 55% | <) And | Ys, | | ☐ Yes XNo Are CT RC | XYes □ No Are MCP A | MA MCP PRESUMPTIVE CE | MA MCP CAM | State /Fed Program Criteria | | | Report Information - Data Deliverables | Date Rec'd in Lab: 🏼 💪 🛮 🕻 |
| N | Date/Time | | | | | | | | | | | | | | | | | Are CT RCP (Reasonable Confidence Protocols) Required? | Are MCP Analytical Methods Required? | ERTAINTY - CT REASON | R66W-1 | Criteria | | l scu | | $G _{V} \cap \mathcal{V}$ alph |
| See raverse side: | | Please print clearly, legibly and com- pletely. Samples can not be legged | | | | | | *moder Sike metal | | | | Sample Specific Comments | (Please specify below) | Preservation Lab to do Preservation Lab to do | Mot poded | SAMPLE HANDLING | | Protocols) Required? | | CERTAINTY - CT REASONABLE CONFIDENCE PROTOCOLS | AND THE STATE OF T | | Clo AP | Same as Client info PO#: | Billing Information | ALPHA JOD #: (100 88) 2 |
| | s are resolved 7 re-subject to 2 | 34 S | | 4 | | | | * | | A Record to the Control of the Contr | W | | mr | | # (| | | | | ROTOCOLS | | | | | | 5812 |

Appendix F
Copy of DEP Transmittal Forms (BWSC-104 and BWSC-108)



BWSC104

RESPONSE ACTION OUTCOME (RAO) STATEMENT

Pursuant to 310 CMR 40.1000 (Subpart J)

| Rele | ease | Tracking | Number |
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For sites with multiple RTNs, enter the Primary RTN above. A. SITE LOCATION: 1. Site Name/Location Aid: _____ 2. Street Address: _ _____ 4. ZIP Code: ___ City/Town: ____ 5. Check here if a Tier Classification Submittal has been provided to DEP for this disposal site. a. Tier IA b. Tier IB c. Tier IC d. Tier II 6. If a Tier I Permit has been issued, provide Permit Number: ___ B. THIS FORM IS BEING USED TO: (check all that apply) 1. List Submittal Date of RAO Statement (if previously submitted): _ mm/dd/yyyy 2. Submit a Response Action Outcome (RAO) Statement a. Check here if this RAO Statement covers additional Release Tracking Numbers (RTNs). RTNs that have been previously linked to a Tier Classified Primary RTN do not need to be listed here. b. Provide additional Release Tracking Number(s) covered by this RAO Statement. 3. Submit a Revised Response Action Outcome Statement a. Check here if this Revised RAO Statement covers additional Release Tracking Numbers (RTNs), not listed on the RAO Statement or previously submitted Revised RAO Statements. RTNs that have been previously linked to a Tier Classified Primary RTN do not need to be listed here. b. Provide additional Release Tracking Number(s) covered by this RAO Statement. 4. Submit a Response Action Outcome Partial (RAO-P) Statement Check above box, if any Response Actions remain to be taken to address conditions associated with this disposal site having the Primary RTN listed in the header section of this transmittal form. This RAO Statement will record only an RAO-Partial Statement for that RTN. A final RAO Statement will need to be submitted that references all RAO-Partial Statements and, if applicable, covers any remaining conditions not covered by the RAO-Partial Statements. Also, specify if you are an Eligible Person or Tenant pursuant to M.G.L. c. 21E s.2, and have no further obligation to conduct response actions on the remaining portion(s) of the disposal site: a. Eligible Person b. Eligible Tenant 5. Submit an optional **Phase I Completion Statement** supporting an RAO Statement 6. Submit a Periodic Review Opinion evaluating the status of a Temporary Solution for a Class C-1 RAO Statement, as specified in 310 CMR 40.1051 (Section F is optional) 7. Submit a Retraction of a previously submitted Response Action Outcome Statement (Sections E & F are not required) (All sections of this transmittal form must be filled out unless otherwise noted above)

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| _ | , v | J | \mathbf{c} | u | - |

| Pursuant to 310 CMR 40.1000 (Subpart S | WE (RAO) STATEMENT | se Tracking Number |
|---|---|--------------------|
| C. DESCRIPTION OF RESPONSE ACTIONS: (check all that 1. Assessment and/or Monitoring Only | t apply; for volumes, list cumulative amounts) 2. Temporary Covers or Caps | |
| 3. Deployment of Absorbent or Containment Materials | 4. Treatment of Water Supplies | , |
| 5. Structure Venting System | 6. Engineered Barrier | , |
| 7. Product or NAPL Recovery | 8. Fencing and Sign Posting | |
| Groundwater Treatment Systems | 10. Soil Vapor Extraction | |
| 11. Bioremediation | 12. Air Sparging | |
| 13. Monitored Natural Attenuation | 14. In-situ Chemical Oxidation | |
| | | |
| 15. Removal of Contaminated Soils | Tatimated valume in aukie varde | |
| a. Re-use, Recycling or Treatment i. On Site E | Estimated volume in cubic yards | |
| ii. Off Site | Estimated volume in cubic yards | |
| iia. Facility Name: | Town: | _ State: |
| iib. Facility Name: | Town: | State: |
| iib. I dointy Haino. | | |
| iii. Describe: | | |
| b. Landfill | | |
| i. Cover Estimated volume in cubic yards | | |
| Facility Name: | Town: | State: |
| | | |
| ii. Disposal Estimated volume in cubic yards | | |
| Facility Name: | Town: | _ State: |
| Tuomity Hamo. | | |
| 16. Removal of Drums, Tanks or Containers: | | |
| a. Describe Quantity and Amount: | | |
| | | |
| | | |
| b. Facility Name: | Town: | _ State: |
| a Facility Names | Town | Stato |
| c. Facility Name: | TOWIT. | _ Olale |
| 17. Removal of Other Contaminated Media: | | |
| a. Specify Type and Volume: | | |
| a. opoony typo and volume. | | |
| | | |
| b. Facility Name: | Town: | State: |
| c. Facility Name: | Town: | State: |

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RESPONSE ACTION OUTCOME (RAO) STATEMENT

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Release Tracking Number

| Pursuant to 310 CMR 40.1000 (Subpart J) | |
|--|--|
| C. DESCRIPTION OF RESPONSE ACTIONS (cont.): (check all that apply; for volumes, list cumulative amounts) | |
| 18. Other Response Actions: | |
| Describe: | |
| 19. Use of Innovative Technologies: | |
| Describe: | |
| D. SITE USE: | |
| 1. Are the response actions that are the subject of this submittal associated with the <i>redevelopment</i> , <i>reuse</i> or the <i>major</i> expansion of the current use of property(ies) impacted by the presence of oil and/or hazardous materials? | |
| a. Yes b. No c. Don't know | |
| 2. Is the property a vacant or under-utilized commercial or industrial property ("a brownfield property")? | |
| a. Yes b. No c. Don't know | |
| 3. Will funds from a state or federal brownfield incentive program be used on one or more of the property(ies) within the disposal site? | |
| a. Yes b. No c. Don't know If Yes, identify program(s): | |
| 4. Has a Covenant Not to Sue been obtained or sought? | |
| a. Yes b. No c. Don't know | |
| 5. Check all applicable categories that apply to the person making this submittal: | |
| b. Community Development Corporation c. Economic Development and Industrial Corporation | |
| d. Private Developer e. Fiduciary f. Secured Lender g. Municipality | |
| h. Potential Buyer (non-owner) i. Other, describe: | |
| | |
| This data will be used by MassDEP for information purposes only, and does not represent or create any legal commitment, obligation or liability on the part of the party or person providing this data to MassDEP. | |
| E. RESPONSE ACTION OUTCOME CLASS: | |
| Specify the Class of Response Action Outcome that applies to the disposal site, or site of the Threat of Release. Select ONLY one Class. | |
| 1. Class A-1 RAO: Specify one of the following: | |
| a. Contamination has been reduced to background levels. b. A Threat of Release has been eliminated. | |
| 2. Class A-2 RAO: You MUST provide justification that reducing contamination to or approaching background levels is infeasible. | |
| 3. Class A-3 RAO: You MUST provide an implemented Activity and Use Limitation (AUL) and justification that reducing contamination to or approaching background levels is infeasible. | |
| 4. Class A-4 RAO: You MUST provide an implemented AUL, justification that reducing contamination to or approaching background levels is infeasible, and justification that reducing contamination to less than Upper Concentration Limits (UCLs) 15 feet below ground surface or below an Engineered Barrier is infeasible. If the Permanent Solution relies upon an Engineered Barrier, you must provide or have previously provided a Phase III Remedial Action Plan that justifies the selection of the Engineered Barrier. | |

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| BWS | C1 | 04 |
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| Pursuant to 310 CMR 40.1000 (Subpart J) |
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| E. RESPONSE ACTION OUTCOME CLASS (cont.): |
| 5. Class B-1 RAO: Specify one of the following: |
| a. Contamination is consistent with background levels b. Contamination is NOT consistent with background levels. |
| 6. Class B-2 RAO: You MUST provide an implemented AUL. |
| 7. Class B-3 RAO: You MUST provide an implemented AUL and justification that reducing contamination to less than Upper Concentration Limits (UCLs) 15 feet below ground surface is infeasible. |
| 8. Class C-1 RAO: You must submit a plan as specified at 310 CMR 40.0861(2)(h). Indicate type of ongoing response actions. |
| a. Active Remedial System b. Active Remedial Monitoring Program c. None |
| d. Other Specify: |
| 9. Class C-2 RAO: You must hold a valid Tier I Permit or Tier II Classification to continue response actions toward a Permanent Solution. |
| F. RESPONSE ACTION OUTCOME INFORMATION: |
| 1. Specify the Risk Characterization Method(s) used to achieve the RAO described above: |
| a. Method 1 b. Method 2 c. Method 3 |
| d. Method Not Applicable-Contamination reduced to or consistent with background, or Threat of Release abated |
| 2. Specify all Soil Category(ies) applicable. More than one Soil Category may apply at a Site. Be sure to check off all APPLICABLE categories: |
| a. S-1/GW-1 d. S-2/GW-1 g. S-3/GW-1 |
| ☐ b. S-1/GW-2 ☐ e. S-2/GW-2 ☐ h. S-3/GW-2 |
| c. S-1/GW-3 f. S-2/GW-3 i. S-3/GW-3 |
| 3. Specify all Groundwater Category(ies) impacted. A site may impact more than one Groundwater Category. Be sure to check off all IMPACTED categories: |
| a. GW-1 b. GW-2 c. GW-3 d. No Groundwater Impacted |
| 4. Specify remediation conducted: |
| a. Check here if soil remediation was conducted. |
| b. Check here if groundwater remediation was conducted. |
| 5. Specify whether the analytical data used to support the Response Action Outcome was generated pursuant to the Department's Compendium of Analytical Methods (CAM) and 310 CMR 40.1056: |
| a. CAM used to support all analytical data. b. CAM used to support some of the analytical data. |
| c. CAM not used. |
| 6. Check here to certify that the Class A, B or C Response Action Outcome includes a Data Usability Assessment and Data Representativeness Evaluation pursuant to 310 CMR 40.1056. |
| 7. Estimate the number of acres this RAO Statement applies to: |

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| BWSC104 | |
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Release Tracking Number RESPONSE ACTION OUTCOME (RAO) STATEMENT Pursuant to 310 CMR 40.1000 (Subpart J) G. LSP SIGNATURE AND STAMP: I attest under the pains and penalties of perjury that I have personally examined and am familiar with this transmittal form, including any and all documents accompanying this submittal. In my professional opinion and judgment based upon application of (i) the standard of care in 309 CMR 4.02(1), (ii) the applicable provisions of 309 CMR 4.02(2) and (3), and 309 CMR4.03(2), and (iii) the provisions of 309 CMR 4.03(3), to the best of my knowledge, information and belief, > if Section B indicates that either an RAO Statement, Phase I Completion Statement and/or Periodic Review Opinion is being provided, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed and implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40,0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal. I am aware that significant penalties may result, including, but not limited to, possible fines and imprisonment, if I submit information which I know to be false, inaccurate or materially incomplete. 1. LSP #: _____ 3. Last Name: _____ First Name: _____ 4. Telephone: ______ 5. Ext.: _____ 6. FAX: ____ 7. Signature: _____ 8. Date: ____ 9. LSP Stamp: mm/dd/vvvv H. PERSON MAKING SUBMITTAL: c. change in the person 1. Check all that apply: a. change in contact name b. change of address undertaking response actions Name of Organization: ______ 3. Contact First Name: ______ 4. Last Name: _____ 6. Title: _____ Street: _______ 8. State: _____ 9. ZIP Code: __ 7. City/Town: ___ 10. Telephone: ______ 12. FAX: _____ 12. FAX: _____

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| K | 1 | RESPONSE ACTION OUTCOME (RAO) STATEMENT | Release Tracking Number |
|-------|---------------|--|-----------------------------------|
| | -(d): | Pursuant to 310 CMR 40.1000 (Subpart J) | |
| . RE | LATIONSHIP 1 | TO RELEASE OR THREAT OF RELEASE OF PERSON MAKING SUBMITTAL: | |
| | 1. RP or PRF | P a. Owner b. Operator c. Generator d. Transpo | rter |
| | | e. Other RP or PRP Specify: | |
| | 2. Fiduciary, | Secured Lender or Municipality with Exempt Status (as defined by M.G.L. c. 21E, | s. 2) |
| | 3. Agency or | Public Utility on a Right of Way (as defined by M.G.L. c. 21E, s. 5(j)) | |
| | 4. Any Other | Person Making Submittal Specify Relationship: | |
| | | | |
| J. RI | EQUIRED ATT | ACHMENT AND SUBMITTALS: | |
| | | re if the Response Action(s) on which this opinion is based, if any, are (were) subval(s) issued by DEP or EPA. If the box is checked, you MUST attach a statement ereof. | |
| | | re to certify that the Chief Municipal Officer and the Local Board of Health have be ement that relies on the public way/rail right-of-way exemption from the requireme | |
| | | re to certify that the Chief Municipal Officer and the Local Board of Health have be ent with instructions on how to obtain a full copy of the report. | en notified of the submittal of a |
| | the Disposal | re to certify that documentation is attached specifying the location of the Site, or the Site subject to this RAO Statement. If submitting an RAO Statement for a PORTI ent the location and boundaries for both the portion subject to this submittal and, e. | ON of a Disposal Site, you |
| | disposal site | e to certify that, pursuant to 310 CMR 40.1406, notice was provided to the owner (boundaries, or notice was not required because the disposal site boundaries are ducting response actions. (check all that apply) | |
| | a. Notice | e was provided prior to, or concurrent with the submittal of a Phase II Completion | Statement to the Department. |
| | b. Notice | e was provided prior to, or concurrent with the submittal of this RAO Statement to | the Department. |
| | C. Notice | e not required. d. Total number of property owners notified, if applicable: | |
| | copy of each | re if required to submit one or more AULs. You must submit an AUL Transmittal implemented AUL related to this RAO Statement. Specify the type of AUL(s) below, B-3 RAO Statements) | |
| | a. Notice | e of Activity and Use Limitation b. Number of Notices submitted: | |
| | c. Grant | of Environmental Restriction d. Number of Grants submitted: | |
| | | Compliance Fee is required for any of the RTNs listed on this transmittal form, chee was submitted to DEP, P. O. Box 4062, Boston, MA 02211. | neck here to certify that an RAO |
| | | re if any non-updatable information provided on this form is incorrect, e.g. Site Aco the DEP Regional Office. | Idress/Location Aid. Send |
| | 9. Check her | re to certify that the LSP Opinion containing the material facts, data, and other info | ormation is attached. |

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| | ISE ACTION OUTCOME (RAO 310 CMR 40.1000 (Subpart J) |) STATEMENT | Release Tracking Number |
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| . CERTIFICATION OF PERSON | MAKING SUBMITTAL: | | |
| xamined and am familiar with the ansmittal form, (ii) that, based of atterial information contained in the I am fully authorized to make thit on whose behalf this submitty on whose behalf this submitty. | , attest under the pa ne information contained in this submitta on my inquiry of those individuals immedi in this submittal is, to the best of my know this attestation on behalf of the entity leg nittal is made am/is aware that there are t, for willfully submitting false, inaccurate | I, including any and all doc ately responsible for obtai ledge and belief, true, acc pally responsible for this su significant penalties, inclu | cuments accompanying this ining the information, the urate and complete, and (iii) ubmittal. I/the person or ding, but not limited to, |
| Ву: | Signature | 3. Title: | |
| For:(Name of person | or entity recorded in Section H) | 5. Date: | mm/dd/yyyy |
| _ | s of the person providing certification is d | ifferent from address reco | rded in Section H. |
| | 9. 9 | State: 10. 2 | ZIP Code: |
| 1. Telephone: | 12. Ext.: | 13. FAX: | |
| BILLABLE YE SECTIONS C | IBJECT TO AN ANNUAL COMPLIANCE AS EAR FOR THIS DISPOSAL SITE. YOU MUS IF THIS FORM OR DEP MAY RETURN THE OMPLETE FORM, YOU MAY BE PENALIZE | T LEGIBLY COMPLETE AL DOCUMENT AS INCOMPL | L RELEVANT ETE. IF YOU |
| Date Stamp (DEP USE ON | JLY:) | | |
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BWSC108

| COMPREHENSIVE RESPONSE ACTION TRANSMITTAL |
|---|
| FORM & PHASE I COMPLETION STATEMENT |

Release Tracking Number

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

| | | LOCATION: |
|-------|--------|---|
| 1. S | Site I | Name: |
| 2. S | Stree | et Address: |
| 3. C | City/ | Town: 4. ZIP Code: |
| | 5 | Check here if a Tier Classification Submittal has been provided to DEP for this disposal site. |
| | J. | |
| | | a. Tier IA b. Tier IB c. Tier IC d. Tier II |
| 6. If | app | olicable, provide the Permit Number: |
| В. Т | HIS | FORM IS BEING USED (check all that apply) |
| | | Submit a Phase I Completion Statement , pursuant to 310 CMR 40.0484. |
| | | Submit a Revised Phase I Completion Statement, pursuant to 310 CMR 40.0484. |
| | | Submit a Phase II Scope of Work, pursuant to 310 CMR 40.0834. |
| | | Submit an interim Phase II Report . This report does not satisfy the response action deadline requirements in |
| | | 0 CMR 40.0500. |
| | 5. | Submit a final Phase II Report and Completion Statement, pursuant to 310 CMR 40.0836. |
| | 6. | Submit a Revised Phase II Report and Completion Statement, pursuant to 310 CMR 40.0836. |
| | 7. | Submit a Phase III Remedial Action Plan and Completion Statement, pursuant to 310 CMR 40.0862. |
| | 8. | Submit a Revised Phase III Remedial Action Plan and Completion Statement, pursuant to 310 CMR 40.0862. |
| | 9. | Submit a Phase IV Remedy Implementation Plan, pursuant to 310 CMR 40.0874. |
| | 10 | Submit a Modified Phase IV Remedy Implementation Plan, pursuant to 310 CMR 40.0874. |
| | 11 | Submit an As-Built Construction Report, pursuant to 310 CMR 40.0875. |
| | 12 | Submit a Phase IV Status Report , pursuant to 310 CMR 40.0877. |
| | 13 | Submit a Phase IV Completion Statement, pursuant to 310 CMR 40.0878 and 40.0879. |
| | | Specify the outcome of Phase IV activities: (check one) |
| | | a. Phase V Operation, Maintenance or Monitoring of the Comprehensive Remedial Action is necessary to achieve a Response Action Outcome. |
| | | b. The requirements of a Class A Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP. |
| | | c. The requirements of a Class C Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) has been or will be submitted to DEP. |
| | | d. The requirements of a Class C Response Action Outcome have been met. Further Operation, Maintenance or Monitoring of the remedial action is necessary to ensure that conditions are maintained and that further progress is made toward a Permanent Solution. A completed Response Action Outcome Statement and Report (BWSC104) has been or will be submitted to DEP. |



COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

| Release | Tracking Number |
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BWSC108

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

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| . THIS | FORM IS BEING USED TO (cont.):(check all that apply) |
| 14 | . Submit a Revised Phase IV Completion Statement, pursuant to 310 CMR 40.0878 and 40.0879. |
| 15 | . Submit a Phase V Status Report , pursuant to 310 CMR 40.0892. |
| 16 | . Submit a Remedial Monitoring Report. (This report can only be submitted through eDEP.) |
| a. | Type of Report: (check one) i. Initial Report ii. Interim Report iii. Final Report |
| b. | Frequency of Submittal: (check all that apply) i. A Remedial Monitoring Report(s) submitted monthly to address an Imminent Hazard. |
| | ii. A Remedial Monitoring Report(s) submitted monthly to address a Condition of Substantial Release Migration.iii. A Remedial Monitoring Report(s) submitted concurrent with a Status Report. |
| C. | Status of Site: (check one) 🔲 i. Phase IV 🔲 ii. Phase V 🔲 iii. Remedy Operation Status 🔲 iv. Class C RAO |
| d. | Number of Remedial Systems and/or Monitoring Programs: |
| | separate BWSC108A, CRA Remedial Monitoring Report, must be filled out for each Remedial System and/or Monitoring rogram addressed by this transmittal form. |
| 17 | . Submit a Remedy Operation Status , pursuant to 310 CMR 40.0893. |
| 18 | . Submit a Status Report to maintain a Remedy Operation Status , pursuant to 310 CMR 40.0893(2). |
| | . Submit a Transfer and/or a Modification of Persons Maintaining a Remedy Operation Status (ROS) , pursuant to 310 MR 40.0893(5) (check one, or both, if applicable). |
| | a. Submit a Transfer of Persons Maintaining an ROS (the transferee should be the person listed in Section D, "Person Undertaking Response Actions"). |
| | b. Submit a Modification of Persons Maintaining an ROS (the primary representative should be the person listed in Section D, "Person Undertaking Response Actions"). |
| C. | . Number of Persons Maintaining an ROS not including the primary representative: |
| 20 | . Submit a Termination of a Remedy Operation Status , pursuant to 310 CMR 40.0893(6).(check one) |
| | a. Submit a notice indicating ROS performance standards have not been met. A plan and timetable pursuant to 310 CMR 40.0893(6)(b) for resuming the ROS are attached. |
| | b. Submit a notice of Termination of ROS. |
| 21 | . Submit a Phase V Completion Statement, pursuant to 310 CMR 40.0894. |
| Sp | ecify the outcome of Phase V activities: (check one) |
| | a. The requirements of a Class A Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement (BWSC104) will be submitted to DEP. |
| | b. The requirements of a Class C Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP. |
| | c. The requirements of a Class C Response Action Outcome have been met. Further Operation, Maintenance or Monitoring of the remedial action is necessary to ensure that conditions are maintained and/or that further progress is made toward a Permanent Solution. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP. |
| 22 | . Submit a Revised Phase V Completion Statement, pursuant to 310 CMR 40.0894. |
| 23 | . Submit a Post-Class C Response Action Outcome Status Report, pursuant to 310 CMR 40.0898. |

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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

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BWSC108

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

| C | I SP | SIGN | ΙΔΤΙ | IRF | | STAMP | ١. |
|----|------|------|-------|-----|------|---------|----|
| v. | LUE | JIGI | יו חי | | AIND | SIAIVIE | |

I attest under the pains and penalties of perjury that I have personally examined and am familiar with this transmittal form, including any and all documents accompanying this submittal. In my professional opinion and judgment based upon application of (i) the standard of care in 309 CMR 4.02(1), (ii) the applicable provisions of 309 CMR 4.02(2) and (3), and 309 CMR 4.03(2), and (iii) the provisions of 309 CMR 4.03(3), to the best of my knowledge, information and belief,

- > if Section B indicates that a **Phase I, Phase II, Phase III, Phase IV or Phase V Completion Statement** and/or a **Termination of a Remedy Operation Status** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed and implemented in accordance with the applicable provisions of M.G.L c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B indicates that a **Phase II Scope of Work** or a **Phase IV Remedy Implementation Plan** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B indicates that an As-Built Construction Report, a Remedy Operation Status, a Phase IV, Phase V or Post-Class C RAO Status Report, a Status Report to Maintain a Remedy Operation Status, a Transfer or Modification of Persons Maintaining a Remedy Operation Status and/or a Remedial Monitoring Report is being submitted, the response action(s) that is (are) the subject of this submittal (i) is (are) being implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal.

I am aware that significant penalties may result, including, but not limited to, possible fines and imprisonment, if I submit information which I know to be false, inaccurate or materially incomplete.

| 1. LSP #: | | |
|----------------------|------------------|--|
| 2. First Name: | 3. Last Name: | |
| 4. Telephone: | 5. Ext.: 6. FAX: | |
| 7. Signature: | | |
| 8. Date:(mm/dd/yyyy) | 9. LSP Stamp: | |
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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

| Release | Tracking | Numbe |
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BWSC108

| Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H) | | | | |
|---|--|--|--|--|
| D. PERSON UNDERTAKING RESPONSE ACTIONS: | | | | |
| 1. Check all that apply: a. change in contact name b. change of address c. change in the person undertaking response actions | | | | |
| 2. Name of Organization: | | | | |
| 3. Contact First Name: 4. Last Name: | | | | |
| 5. Street: 6. Title: | | | | |
| 7. City/Town: 9. ZIP Code: | | | | |
| 10. Telephone: 11. Ext.: 12. FAX: | | | | |
| E. RELATIONSHIP TO SITE OF PERSON UNDERTAKING RESPONSE ACTIONS: Check here to change relationship | | | | |
| 1. RP or PRP a. Owner b. Operator c. Generator d. Transporter | | | | |
| e. Other RP or PRP Specify: | | | | |
| 2. Fiduciary, Secured Lender or Municipality with Exempt Status (as defined by M.G.L. c. 21E, s. 2) | | | | |
| 3. Agency or Public Utility on a Right of Way (as defined by M.G.L. c. 21E, s. 5(j)) | | | | |
| 4. Any Other Person Undertaking Response Actions Specify Relationship: | | | | |
| F. REQUIRED ATTACHMENT AND SUBMITTALS: | | | | |
| Check here if the Response Action(s) on which this opinion is based, if any, are (were) subject to any order(s), permit(s) and/or approval(s) issued by DEP or EPA. If the box is checked, you MUST attach a statement identifying the applicable provisions thereof. | | | | |
| 2. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the submittal of any Phase Reports to DEP. | | | | |
| 3. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the availability of a Phase III Remedial Action Plan. | | | | |
| 4. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the availability of a Phase IV Remedy Implementation Plan. | | | | |
| 5. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of any field work involving the implementation of a Phase IV Remedial Action. | | | | |
| 6. If submitting a Transfer of a Remedy Operation Status (as per 310 CMR 40.0893(5)), check here to certify that a statement detailing the compliance history for the person making this submittal (transferee) is attached. | | | | |
| 7. If submitting a Modification of a Remedy Operation Status (as per 310 CMR 40.0893(5)), check here to certify that a statement detailing the compliance history for each new person making this submittal is attached. | | | | |
| 8. Check here if any non-updatable information provided on this form is incorrect, e.g. Site Name. Send corrections to: BWSC.eDEP@state.ma.us. | | | | |
| 9. Check here to certify that the LSP Opinion containing the material facts, data, and other information is attached. | | | | |
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Release Tracking Number

COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

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| | .0484 (Subpart D) and 40.08 | · , | | |
| G. CERTIFICATION OF PERSON UNDERT | | | | |
| accompanying this transmittal form, (ii) that, information, the material information contains complete, and (iii) that I am fully authorized t | based on my inquiry of those ed in this submittal is, to the o make this attestation on be tal is made am/is aware that | nder the pains and penalties of perjury (i) that I have nis submittal, including any and all documents se individuals immediately responsible for obtaining the elbest of my knowledge and belief, true, accurate and behalf of the entity legally responsible for this submittal. I/the t there are significant penalties, including, but not limited to, e, or incomplete information. | | |
| perjury that I am fully authorized to act on be | half of all persons performing espondence from MassDEP | ion Status (ROS), I attest under the pains and penalties of ng response actions under the ROS as stated in 310 CMR with respect to performance of response actions under the | | |
| | , and I am aware that there a | rom MassDEP shall be deemed received by all the persons are significant penalties, including, but not limited to, possible aplete information. | | |
| 2. By: | | 3. Title: | | |
| • | nature | | | |
| 4. For: | | 5. Date: | | |
| (Name of person or entity recorded in Section D) | | (mm/dd/yyyy) | | |
| 7. Street: | | 10 710 0 | | |
| 8. City/Town: | 9. | 0. State: 10. ZIP Code: | | |
| 11. Telephone: | 12. Ext.: | 13. FAX: | | |
| YOU ARE SUBJECT TO AN ANNUAL COMPLIANCE ASSURANCE FEE OF UP TO \$10,000 PER BILLABLE YEAR FOR THIS DISPOSAL SITE. YOU MUST LEGIBLY COMPLETE ALL RELEVANT SECTIONS OF THIS FORM OR DEP MAY RETURN THE DOCUMENT AS INCOMPLETE. IF YOU SUBMIT AN INCOMPLETE FORM, YOU MAY BE PENALIZED FOR MISSING A REQUIRED DEADLINE. | | | | |
| Date Stamp (DEP USE ONLY:) | | | | |
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