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Phase II Comprehensive Site Assessment Addendum for Demolition Debris Area

RTN 4-3024222 Former Bird Machine Company Site Walpole, MA

Submitted to:

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LIST OF ACRONYMS

ACM asbestos containing material

AIS asbestos in soil

AMEC AMEC Environment and Infrastructure

AUL Activity and Use Limitation bgs below ground surface BHI Baker Hughes Inc.
BMC Bird Machine Company

CAM Compendium of Analytical Methods

CIH Certified Industrial Hygienist

CMR Code of Massachusetts Regulations CSA Comprehensive Site Assessment

CSM Conceptual Site Model
COC contaminant of concern
DDA Demolition Debris Area
DUA data usability assessment

EPH extractable petroleum hydrocarbon

ft feet

IRA Immediate Response Action

LRA Lead Release Area

MADEP Massachusetts Department of Environmental Protection

MBA Manufacturing Building Area
MCP Massachusetts Contingency Plan
MS/MSD matrix spike/matrix spike duplicate

NSR No Significant Risk

PAH polycyclic aromatic hydrocarbon

PCB polychlorinated biphenyl
PLM polarized light microscopy
RAM Release Abatement Measure
RAO Response Action Outcome
RC Risk Characterization

REDUA representativeness and data usability assessment

RTN Release Tracking Number SDG sample delivery group SRS South Rail Spur

STL Severn Trent Laboratories
SVOC semivolatile organic compound
TEM transmission electron microscopy

U.S. EPA United States Environmental Protection Agency

VOC volatile organic compound VPH volatile petroleum hydrocarbon

Weston Weston Solutions, Inc.



EXECUTIVE SUMMARY

On behalf of Baker Hughes, Inc. (BHI), AMEC Earth and Environmental, Inc. (AMEC) has completed a Phase II Comprehensive Site Assessment (CSA) Addendum for the portion of the former Bird Machine Company (BMC) Property located in Walpole, Massachusetts known as the Demolition Debris Area (DDA). The DDA is an exposure area and is a portion of the site assigned Release Tracking Number (RTN) 4-3024222 under the Massachusetts Contingency Plan (MCP). This Phase II CSA Addendum serves as an update to a July 2007 Phase II CSA (Weston 2007). It presents the data collected at the DDA from June 2007 to the present, updates to the Phase II CSA as a result of the additional data collected, and an updated risk characterization. A Phase II CSA addressing three other exposure areas, the manufacturing building area (MBA), the lead release area 3 (LRA3), and the south rail spur (SRS) was finalized on October 18, 2011. These two Phase II CSAs together characterize the "Site", which is represented by the single unclosed RTN (4-3024222) at the property.

The Phase II CSA Addendum addresses volatile organic constituents (VOCs), semi-volatile organic constituents (SVOCs), extractable petroleum hydrocarbons (EPH), polycyclic aromatic hydrocarbons (PAHs), dioxin/furan congeners, and various metals detected in soil and groundwater samples collected from the DDA. The Phase II CSA also includes evaluations of asbestos in soil (AIS) identified within this exposure area.

Data from site investigations completed by AMEC, site assessment activities completed by Weston Solutions, Inc. of Concord, New Hampshire (Weston), and information from other sources (e.g., Massachusetts Department of Environmental Protection [MADEP] and United States Environmental Protection Agency [U.S. EPA] guidance documents), were used to complete the CSA.

In accordance with the requirements of 310 CMR 40.0000 Subpart I of the MCP, a Method 3 risk characterization (RC) of harm to human health, public welfare, safety, and the environment was completed. This RC replaces a Method 1 RC that was prepared by Weston. The Method 1 RC addressed soil and groundwater at the DDA and was not able to conclude NSR. Additionally, Weston determined that a Method 3 RC would eventually be necessary to support a response action outcome (RAO) statement due to the presence of dioxins (which are bioaccumulative) in the top two feet of soil, and the presence of asbestos. This Method 3 RC has been conducted assuming that an Activity and Use Limitation (AUL) will be implemented at the DDA prohibiting disruption of the ground surface. The risk characterization therefore does not evaluate any Site use or development other than incidental trespassing.

The Phase II investigations presented in this Phase II CSA Addendum consisted of ground water sampling in and around the DDA and soil sampling for asbestos within the DDA.

The nature and extent of soil contamination at the DDA was previously documented in Weston's 2007 Phase II CSA. Additional asbestos sampling has demonstrated that asbestos could be present within the footprint of the DDA fill area both horizontally and vertically. Groundwater sampling results are consistent with Weston's characterization in the 2007 Phase II CSA. Groundwater contains metals and sporadic PAH detections. Weston's 2007 Phase II CSA determined that these groundwater contaminants do not represent a significant source of contamination to river sediment or surface water.



The results of the human health and environmental risk characterization indicate that a condition of No Significant Risk (NSR) of harm to health, safety, public welfare, and the environment has been achieved at the DDA.



1.0 INTRODUCTION

On behalf of BHI, AMEC has completed a Phase II CSA Addendum pursuant to 310 Code of Massachusetts Regulations (CMR) 40.0830 of the Massachusetts Contingency Plan (MCP). This Phase II CSA Addendum serves as an update to a July 2007 Phase II CSA for the DDA portion of RTN 4-3024222 (Weston 2007). It presents the data collected at the DDA from June 2007 to the present, updates to the 2007 Phase II CSA as a result of the additional data collected, and an updated risk characterization. The property location is provided in Figure 1-1.

At the time of the July 2007 CSA, DDA was assigned RTN 4-3024105 and was classified as a Tier II Disposal Site under the MCP. The DDA was linked to RTN 4-3024222 in the January 2008 Tier IB Permit Application for the Site. A Phase II CSA addressing other exposure areas in this linked RTN – the manufacturing building area (MBA), the lead release area (LRA), the south rail spur (SRS), and the Neponset River – was finalized on October 18, 2011. These two Phase II CSAs together characterize RTN 4-3024222 at the property. Figure 1-2 presents all five exposure areas described in these two CSAs.

This Phase II CSA Addendum documents field investigations completed between June 2007 and the present and the resulting findings and laboratory analytical results. Any information collected prior to June 2007 for the DDA was presented in Weston's 2007 Phase II CSA. This report follows the same outline as Weston's 2007 report and only those sections where additional information was available were updated. Those sections that did not change indicate that there was no update from the 2007 Phase II CSA.



2.0 GENERAL SITE INFORMATION [310 CMR 40.0835(4)(A)]

There are no updates to Section 2.0 of Weston's 2007 Phase II CSA for the DDA.



3.0 SITE HISTORY [310 CMR 40.0835(4)(C)]

There are no updates to Section 3.0 of Weston's 2007 Phase II CSA for the DDA.



4.0 INVESTIGATIONS [310 CMR 40.0835(4)(D)1]

Weston's 2007 Phase II CSA for the DDA presents a discussion of the sampling activities conducted at the DDA through June 2007. The following sections present additional sampling that was conducted at the DDA after June 2007.

4.1 Investigation Activities and Results

4.1.1 Asbestos Sampling – April 2011

The Phase II CSA prepared by Weston in July 2007 was unable to conclude No Significant Risk (NSR) due to, among other things, the visually observed presence of asbestos in soil. In order to address the presence of asbestos at the Site and evaluate the risks associated with its presence, AMEC performed Asbestos in Soil (AIS) Investigations at the Site. During the week of April 25th, 2011, Certified Industrial Hygienist (CIH) / Certified Safety Professional (CSP) Michael Matilainen, of AMEC performed asbestos sampling assisted by Environmental Scientist, Alexander Ranieri.

AMEC collected soil samples at 42 locations, 20 of which were located at or just beyond the defined DDA perimeter, and 22 of which were collected within the DDA areas (further defined as the Eastern Clearing, Central Clearing, and Western Clearing areas). Figure 4-1 presents these soil sampling locations. Sample locations were predetermined based on a grid pattern and located in the field using a global positioning system (GPS) device. Soil samples were collected via hand excavation for lab analysis. At each location, samples were collected from two depth intervals, 0-3 inches (interval "A") and 3-24 inches (interval "B"). Note that additional sample was collected from locations 8 (both A & B) and 14 (both A & B). These samples were originally intended for Transmission Electron Microscopy (TEM) TEM analysis, and as such have "-TEM" at the end of each sample ID. However, this "-TEM" suffix does not necessarily correspond to TEM analysis, as some of these samples were submitted for Polarized Light Microscopy (PLM) analysis only. Sample duplicates were collected at a rate of approximately ten percent resulting in twelve duplicates from the following locations: 1A, 2A, 3A, 4A, 5A, 6A, 7A, 8A-TEM, 9A, 38A, 41B, and 42A. Due to the visual presence of asbestos containing material (ACM) at sampling Location 8, a surface sample and duplicate surface sample were collected from the leaf litter (samples 8AS- AR and 8AS-TEM) on the soil surface (above 0 inches).

These selected depth intervals were considered appropriate since the property is not proposed for any future intrusive activities or disturbance (such as construction or grading); an Activity and Use Limitation (AUL) to that effect will be implemented as part of the final Response Action Outcome (RAO) Statement. Samples from 0-24 inches adequately characterize surficial material that may potentially be contacted or suspended due to disturbance by trespassers or other transient Site occupants. Samples from 0-3 inches specifically address the potential for airborne asbestos to present an inhalation hazard under current and future Site conditions.

Samples were shipped to International Asbestos Testing Laboratories (IATL) of Mount Laurel, New Jersey, an asbestos-accredited laboratory. All samples were initially analyzed for total asbestos using PLM U.S. EPA Method 600/R-93/116. Mr. Matilainen, in conjunction with recommendation from IATL Laboratories, then selected an additional seven samples (Samples



8A-TEM, 8A-TEM-DUP, 8B-TEM, 14A-TEM, 14B-TEM, 33A, and 33B) for more sensitive analysis and specific asbestos type characterization via TEM. Samples were analyzed by an IATL proprietary TEM soil method based upon U.S. EPA Region 1 Asbestos in Soils, Sludge and Sediments, 1994. Table 4-1 presents a summary of soil samples and analyses¹.

PLM and TEM analysis results give the percent and type of asbestos in the sample. The asbestos content of manufactured items ranges from one percent to 100 percent. Materials containing asbestos greater than one percent (>1%) by weight are considered ACM by the U.S. EPA.

The only area where asbestos was detected visually or by either PLM or TEM was at Location 8. Table 4-2 presents a summary of results at this location. As described above, a total of seven samples were collected from Location 8 (two surficial debris samples, three soil samples from 0-3 inches which include one duplicate, and two soil samples from 3-24 inches). Chrysotile asbestos, the most common type of asbestos found in buildings (often defined as "white asbestos") was the primary type of asbestos found in this location. The following is a summary of the asbestos sampling results at location 8.

Asbestos was detected in both surficial samples with the following results: chrysotile asbestos was detected at 1.9% in 8AS-AR and chrysotile asbestos was detected at 1.7% and crocidolite was detected at 5.2% in sample 8AS-TEM (all detected via PLM).

Asbestos was detected in one of the three 0-3 inch samples (8A) at 1.3% as chrysotile via PLM. Trace amounts (<1%) of chrysotile asbestos were detected in the other two samples (8A-TEM and 8A-TEM-DUP) via PLM.

Chrysotile was detected at trace amounts in one of the associated deep samples (8B-TEM) collected from 3-24 inches via PLM. However, chrysotile asbestos was not detected in the other deep sample (8B) via PLM.

Of the seven samples that were submitted for TEM analysis, only two samples detected asbestos via this method. Samples 8B and 8A-TEM-DUP detected trace amounts of chrysotile asbestos via TEM analysis.

4.1.2 Asbestos Elutriator Results

Based on the results of the April 2011 asbestos sampling, asbestos was determined to be present in soil. Given the small area of AIS impact, the surficial presence of asbestos at the DDA is unlikely to be an exposure concern, especially given the unoccupied nature of the Site and the nature of the asbestos observed, which consisted of larger pieces (asbestos Transite board and fiber parts around 0.5 square inches in size) that are unlikely to become airborne. However, since asbestos was confirmed at >1% at Location 8, a condition of NSR could not be

¹ Because some submitted samples intended to be held for TEM analysis were analyzed by the laboratory by PLM, the sample IDs are not necessarily indicative of the method used. Table 4-1 clearly identifies the sample IDs and associated analyses.



demonstrated without further testing. AMEC conducted additional analysis that further characterizes the exposure potential associated with asbestos within the DDA.

To address the potential for risk, aliquots of the three DDA soil samples that had asbestos detections of >1% (sample 8A [0-3 inch surface soil] and samples 8AS-AR and 8AS-TEM [surface debris]) were tested for airborne fibers using the elutriator method². Analyses were performed by EMS Laboratories of Pasadena, CA. This method provides "concentration measurements for the specific set of asbestos structure sizes and shapes that contribute to adverse biological effects" and is specifically designed for risk evaluation purposes. Specifically, it uses a dust generator (tumbler) to estimate the asbestos released from a soil sample during disturbance.

The analytical report from EMS Laboratories is presented in Appendix A. For the three samples analyzed, total asbestos fibers >5 micrometers (µm) ranged from 11.7 to 82.9 million fibers per gram. The sample with the maximum asbestos content was 8AS-TEM (surface debris). Table 4-3 presents a summary of the elutriator results for the asbestos analysis.

4.1.3 Groundwater Sampling

Weston's 2007 Phase II CSA for the DDA discussed investigation activities conducted within the DDA beginning in 1985 through June 2007. This Phase II CSA Addendum covers activities conducted at the Site from June 2007 to the present, that were not presented in the earlier CSA report. A total of four rounds of groundwater sampling have been conducted since June 2007. Copies of laboratory data reports are included in Appendix A.

On June 5 and 6, 2007, DD-MW-002, DD-MW-201, DD-MW-203, DD-MW-204, DD-MW-205, DD-MW-206, and DD-MW-207 were sampled for semivolatile organic compounds (SVOCs), extractable petroleum hydrocarbons (EPH), and dissolved metals. Groundwater samples were collected from upgradient (DD-MW-204 and DD-MW-205) and within and/or downgradient (DD-MW-002, DD-MW-201, DD-MW-203, DD-MW-206, and DD-MW-207) of the known fill area, depicted by the original RTN boundary of the DDA on Figure 4-2. The groundwater samples were analyzed by Severn Trent Laboratories, Inc. (STL) of Westfield, MA.

On June 25, 2007, DD-MW-208 was sampled and a duplicate was collected. The samples were submitted for analysis of SVOCs, volatile organic compounds (VOCs), EPH, and dissolved metals. DD-MW-208 was sampled again on July 23, 2007 and analyzed for SVOCs and EPH. The groundwater samples were analyzed by STL of Westfield, MA.

On December 11, 2007, DD-MW-201, DD-MW-203, DD-MW-207, and DD-MW-208 were sampled for SVOCs, EPH (carbon ranges only), and dissolved metals. A duplicate was collected at DD-MW-207. The groundwater samples were analyzed by Test America of Westfield, MA.

² Berman, DW and Kolk, A. Modified Elutriator Method for the Determination of Asbestos in Soils and Bulk Material. Aeolus, Inc., Albany, California and EMS Laboratories, Pasadena, California. Revision 1. http://www.aeolusinc.com/Modified_Elutriator_Method.pdf.



On May 19, 2008, DD-MW-002, DD-MW-201, DD-MW-203, DD-MW-204, DD-MW-205, DD-MW-206, DD-MW-207, and DD-MW-208 were sampled for SVOCs, EPH (carbon ranges only), dissolved metals, and 1,4-dioxane. A duplicate was collected at DD-MW-201 and DD-MW-207. The groundwater samples were analyzed by Test America of Westfield, MA.

SVOCs were analyzed via U.S. EPA Method 8270C, EPH was analyzed via the MassDEP EPH method, dissolved metals were analyzed via U.S. EPA Method 6010, mercury was analyzed via U.S. EPA Method 7470a, VPH was analyzed via the MADEP volatile petroleum hydrocarbon (VPH) method, and VOCs were analyzed via USEPA Method 8260.

Historically, groundwater analytical results have detected polynuclear aromatic hydrocarbons (PAHs), dissolved metals, toluene, and bromoform in DDA groundwater. The most recent results confirm the detections of PAHs, and dissolved metals. Toluene and bromoform were not detected in any of the sampling rounds presented in this Phase II CSA Addendum. Refer to Table 4-4 for the groundwater analytical results presented in this Phase II CSA Addendum. Figure 4-2 presents the monitoring well locations at the DDA.

4.2 Field Methodology

Details regarding field methodology for work conducted prior to 2007 are included in the previously submitted Phase II CSA for the DDA.

The sampling activities associated with this Phase II Addendum were performed in accordance with MassDEP guidance and methods. Soil samples were collected via hand excavation using clean sampling tools and placed directly into sample containers. Groundwater samples were collected from monitoring wells using low-flow sampling techniques in accordance with U.S. EPA guidelines (U.S. EPA, 1996). Soil and groundwater samples were collected with sufficient quality assurance/quality control to meet MCP Presumptive Certainty criteria.



5.0 SITE HYDROGEOLOGICAL CHARACTERISTICS [310 CMR 40.0835(4)(D)3.]

There are no updates to Section 5.0 of Weston's 2007 Phase II CSA for the DDA.



6.0 NATURE AND EXTENT OF CONTAMINATION [310 CMR 40.0835(4)(F)]

6.1 Nature and Extent of Soil Contamination

The nature and extent of soil contamination was discussed in the 2007 Phase II CSA for the DDA for metals, PAHs, VOCs, SVOCs, EPH and VPH, dioxins, and polychlorinated biphenyls (PCBs). As discussed in the Phase II the boundary of the DDA was identified based on topography; clearings observed in historical aerial photographs; test pit and soil boring sampling conducted during investigations, RAMs, and IRAs; and geophysical surveys. This boundary is consistent with the observed presence of an artificial fill layer in this area.

The results of the asbestos delineation investigation identified ACM and asbestos fibers in soil (greater than 1%) in shallow soils [0-3 inches below ground surface (bgs)] in the vicinity of sampling location number 8. Trace amounts of asbestos were identified in samples from 3-24 inches bgs in the vicinity of sampling location number 8. The purpose of this AIS investigation was to delineate the extent of asbestos in the surface at the DDA. With the exception of sampling Location 8, asbestos was not detected or visually observed within sample locations both within and outside of the DDA boundary.

As described in the previous reports prepared by Weston, asbestos and ACM is a known component of the fill at the DDA. Although response actions during an IRA in 2005 removed 1,106 tons of asbestos-containing fill, Weston was not able to remove all of the asbestos-containing fill materials within the boundary of the DDA. The AIS sampling has confirmed that the surficial extent of asbestos and ACM has been delineated, and is most likely concentrated around Location 8; however, the Site has extensive vegetative cover throughout the inspected areas, which prohibited visual inspection of the entire DDA. While thorough inspections were completed in the sample locations, additional asbestos could be hidden in areas under vegetation. Although concentrations of asbestos appear to decrease with depth, asbestos could also be present in soil below 24 inches. AMEC has conservatively assumed that asbestos could be present within the footprint of the DDA fill area (the original RTN boundary) both horizontally and vertically.

6.2 Nature and Extent of Groundwater Contamination

As described in the Phase II CSA for the DDA, groundwater contamination at the DDA consists of dissolved metals, PAHs, and minimal VOCs. Table 6-1 presents a summary of the most recent DDA groundwater quality results.

All groundwater samples from the most recent sampling rounds contained detectable concentrations of one or more metals. In the 2007 Phase II CSA, detected metals included antimony, arsenic, barium, beryllium, cadmium, chromium, nickel, selenium, silver, thallium, vanadium, and zinc. The recent samples included some of these same metals, but not antimony, beryllium, cadmium, silver, and thallium. Two new metals were detected in the recent samples: mercury once at less than 1 part per billion (ppb) in DD-MW-201, and lead once at an estimated concentration below the 1 ppb reporting limit in DD-MW-208. Mercury was not detected in a subsequent sample from DD-MW-201. Recent concentrations of the other detected metals were similar to the levels reported in 2007; less than 10 ppb except for barium



and chromium. Barium (19-47 ppb) and chromium (1.9-35 ppb) levels were still highest in the wells located inside the DDA (DD-MW-001 and -002), but appeared to be declining over time.

A total of 13 SVOCs and two EPH carbon ranges were detected in groundwater samples collected from June 2007 to May 2008. Nine of these analytes (2,4-dinitrophenol, benzo(a)anthracene, benzo(g,h,i)pyrene, butylbenzylphthalate, chrysene, dibenzo(a,h,)anthracene, di-n-butyl phthalate, di-n-octyl phthalate, and indeno(1,2,3) pyrene) were detected only once in groundwater. Eight of the nine (all except 2,4-dinitrophenol) were detected in downgradient monitoring wells DD-MW-207 and DD-MW-208. 2,4-Dinitrophenol was detected in DD-MW-204 in 2008. All of these analytes had a low frequency of detection (less than 10%) and most were only detected once out of all of the groundwater sampling rounds. Bis(2-ethylhexyl)phthalate was detected in six wells (DD-MW-201, DD-MW-203, DD-MW-204, DD-MW-206, DD-MW-207, and DD-MW-208) in 2008. Fluoranthene was detected three times in 2007 using Method 8270 (DD-MW-207, DD-MW-208 and its duplicate); it was also detected in DD-MW-208 and its duplicate via the EPH analysis as well. Phenanthrene, C11-C22 Aromatic EPH, and C19-C36 Aliphatic EPH were detected in DD-MW-208 and its duplicate during the June 2007 sampling round.

As indicated in Weston's 2007 Phase II CSA, eight PAHs were detected in one groundwater sample collected in June 2006 from DD-MW-203, but the results of additional sampling later in the year suggested that the June 2006 results were an anomaly. The sporadic detections of PAHs in 2007-2008 suggest that 2006 results may not have been anomalous and low levels of PAHs may be present in groundwater at the Site. However, the infrequent and low detections of PAHs suggest that there is not a continuous plume of these contaminants at the DDA.

Both bromomethane and toluene were previously detected in groundwater samples in the DDA. Neither of these compounds were detected in the most recent sampling rounds, and as such are not considered to be a concern.

6.3 Nature and Extent of DDA Contamination in Surface Water/Sediment

As was indicated in Weston's 2007 Phase II CSA, the potential for impacts from the DDA to surface water or sediment has been evaluated and it has been confirmed that there are no impacts to surface water or sediment from the DDA. No new surface water or sediment data were collected since the 2007 CSA.



7.0 ENVIRONMENTAL FATE AND TRANSPORT [310 CMR 40.0835(4)(E)]

A complete discussion of the environmental fate and transport of the contaminants of concern (COCs) at the DDA were discussed in Weston's 2007 Phase II CSA. The additional data that were collected between Weston's 2007 Phase II CSA and this Phase II CSA Addendum did not encounter any new groups of analytes; as such, there are no updates to the Environmental Fate and Transport section.



8.0 CONCEPTUAL SITE MODEL

8.1 Introduction

There are no updates to Section 8.1 of Weston's Phase II 2007 CSA for the DDA.

8.2 Conceptual Site Model

Weston's 2007 Phase II CSA for the DDA presented a Conceptual Site Model (CSM) which is largely confirmed by the soil and groundwater sampling described in this Addendum, which was conducted after June 2007. Soil sampling for asbestos confirmed that this contaminant is present within the DDA boundary. The recent asbestos testing provided additional data on the type of ACM and the potential for airborne fibers as described in Sections 4 and 6.

Groundwater sampling confirmed that metals are the principal COCs, and concentrations appear to be stable or declining. Low-level detections of PAHs in groundwater were previously thought to be anomalous but have continued sporadically, and so are believed to be representative of groundwater. These compounds were included in the earlier CSM primarily as a soil contaminant, though the potential for migration to groundwater was recognized. The infrequent and low detections of PAHs suggest that there is not a continuous plume of these contaminants at the DDA.



9.0 EXPOSURE ASSESSMENT [310 CMR 40.0835(4)(G)]

Section 9.0 of Weston's 2007 Phase II CSA has been combined with Section 10.0 for this Phase II CSA Addendum. Refer to Section 10.0 for a discussion of the exposure assessment.



10.0 RISK CHARACTERIZATION [310 CMR 40.0835(4)(H)]

In accordance with the requirements of 310 CMR 40.0000 Subpart I of the MCP, a Method 3 Risk Characterization (RC) of harm to human health, public welfare, safety, and the environment has been completed. This Method 3 RC is included as Appendix B. A summary of the findings of the RC is presented below.

The Method 3 human health RC of soil and groundwater in the DDA evaluated potential exposures to current and future trespassers, and future hypothetical groundwater use as a potable supply. The results of the human health risk characterization indicate that a condition of No Significant Risk can be demonstrated for soil and groundwater at DDA.

The results of the evaluation of risk of harm to safety and public welfare indicates that no unsafe or nuisance conditions exist at the Site. Soil and groundwater constituent concentrations are less than their respective UCLs. As such, a condition of No Significant Risk to public welfare and a condition of No Significant Risk to safety can be demonstrated at the Site.

The evaluation of potential risk of harm to the environment included a Stage I screening evaluation of the presence of ecological receptors and potential habitat for terrestrial ecological receptors. The Stage I screening demonstrated that potential risk to the environment could not be ruled out, so a Stage II Environmental Risk Characterization (ERC) was performed.

In the Stage II ERC, potential exposures of herbivorous mammals, herbivorous avians, insectivorous mammals, insectivorous avians, carnivorous mammals, and carnivorous avians to COPECs in soil and the food web were evaluated using a hazard quotient approach. Hazard quotients are below 1 for all receptors except the short-tailed shrew. For the shrew, the NOAEL HQs exceeded 1 in DDA for only two COCs, and these HQs were below 10. These exposures are not expected to cause adverse environmental impacts to short-tailed shrew populations or populations at the Site, because the hazard quotient approach used in this evaluation is based on a sensitive individual receptor. Conservative exposure assumptions are combined with conservative toxicity assumptions, so that the resulting risk estimates overestimate potential effects to the populations. HQs within an order of magnitude of 1 are not likely to be associated with population effects. Based on the finding that population-level effects for all receptors are not expected, a condition of No Significant Risk of harm to the environment exists.

This human health and environmental risk characterization concludes that DDA achieves a condition of No Significant Risk.



11.0 DATA USABILITY AND REPRESENTATIVENESS

Data described in this Phase II CSA Addendum and in the original July 2007 Phase II CSA will be used in support of an RAO Statement for portions of the BMC Property. These data were collected after 2003 when the MADEP established "presumptive certainty" requirements as defined in "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data" (BWSC-CAM-VIIA, rev. 3.1 dated May 22, 2003). Copies of lab data packages used to support the RC for the DDA are provided in Appendix C of Weston's 2007 Phase II CSA for data collected up to June 2007, and are included in Appendix A of this Addendum for samples collected after that time.

Per 310 CMR 40.1056(2)(k), a representativeness evaluation and data usability assessment (REDUA) must be conducted for data that are used to support an RAO. MassDEP Policy #WSC-07-350, MCP Representativeness Evaluations and Data Usability Assessments, provides guidance on conducting REDUAs under the MCP. Appendix V of the MassDEP document is a REDUA worksheet.

REDUA has already been conducted for the data presented in Weston's July 2007 Phase II CSA for the DDA, though it was not documented in the REDUA worksheet format published by MassDEP in September 2007. As indicated in Section 9.2.1.1 of the 2007 CSA, Weston used only CAM-compliant, usable, and representative data, which included Weston's data collected from December 2004 through June 2007. Weston determined that data were CAM-compliant by confirming laboratory certification of data packages and by chemist review of each lab report. Weston removed results for individual analytes if they determined the results were not CAM-compliant. Weston also determined that data collected during these investigations were representative of the site conditions, and identified any results that had been affected by removal actions or were otherwise no longer representative of site conditions. Table 4-1 of Weston's 2007 Phase II CSA presents a summary of the soil and groundwater samples collected to support their Phase II CSA.

The information presented in Weston's 2007 Phase II CSA and in this Addendum have been compiled to complete the REDUA provided below. The following subsections provide a representativeness evaluation (11.1) and data usability assessment (11.2) in a format that coincides with the categories specified in the MassDEP worksheet. Appendix C presents the MassDEP worksheet for reference.

11.1 Representativeness Evaluation

A-1 Conceptual Site Model (CSM)

- Placement of fill at the Site or in the vicinity is believed to have started in the late 19th century coincident with the construction of the railroad and industrial development of the Property, and to have ended in the 1970s. The types of materials found in fill in the DDA include fill, demolition debris, machining waste, and testing waste. Some wastes were disposed in containers including 55-gallon drums. The types of contaminants include metals, PAH, EPH, dioxins, and asbestos.
- The fill at the Site was placed directly on the native soil surface. Soil contamination is thus of primary concern. However, other media that would have possibly been impacted



by the fill would include groundwater beneath the Site and surface water and sediment in Cedar Swamp Brook which bounds the Site to the north. Human or ecological receptors would be exposed to contaminants in soil, sediment, or surface water by direct contact or ingestion, and potentially through inhalation of suspended soil (dust).

- The sampling program was designed to characterize the nature and extent of contamination and evaluate potential migration pathways of contamination. Each release mechanism from waste source materials to impacted media was considered and corresponding sampling was conducted to the related media matrix. The results of the sampling program were used to verify and update the CSM. The data supports the following contamination migration pathways: from waste source materials to soil or groundwater, and from contaminated soil to groundwater. The data support the conclusion that contamination migration via groundwater flow or surface runoff, to surface water or sediment, is not a complete pathway.
- Waste source materials in drums and soils were transported off-site during RAM activities beginning in 2005. Approximately 2,191 tons of soil and debris were removed. The excavation ranged from 8-10 feet in depth in the western and eastern clearings and 13 feet in depth in the central clearing. All of the excavations were above the water table. Post excavation, contaminants in soil included metals, dioxin, and petroleum compounds. Concentrations in the depth interval of 0-5 ft bgs were higher than in the depth interval of 5-15 ft bgs.
- Asbestos-Containing Material (ACM) was encountered within the limits of the DDA fill, typically three feet in depth. Approximately 1,106 tons of ACM and soil were transported off-site during an IRA in 2005. Results of the recent asbestos sampling confirm that ACM is still present within the DDA boundary

A-2 Use of Field Screening Data

- The 2011 soil samples were visually screened for the presence of asbestos in addition to laboratory analysis of the samples. The visual observations correlated with the asbestos analytical results in terms of identifying a single location where asbestos was present.
- Weston performed field screening and visual observation of soil to select samples for laboratory analysis, during test pitting and soil boring advancement between 2004 and 2006. Some of the selected samples contained detectable levels of VOCs, SVOCs, dioxins, and metals based on laboratory analysis. Field screening was not used for characterizing exposure point concentrations.

A-3 Sampling Rationale in Support of RAO

- The boundary of the DDA was initially defined based on topography and clearings observed in historical air photos indicating potential disposal activities. The boundary that was initially developed based on these data was further supported by subsurface investigations confirming the presence of an artificial fill layer inside the DDA boundary.
- The analytical results of soil samples collected from inside the DDA boundary exhibited distinct characteristics, including elevated concentrations of metals, PAH/EPH, and dioxins as compared to samples collected from outside the DDA. RAM and IRA activities were conducted to remove buried waste barrels and significant quantities of source materials from the Site. Post-excavation soil samples indicate a significant reduction of the residual concentrations of these contaminants.



- The sampled area around the DDA boundary is called the Perimetric Area. All
 contaminants detected in soil samples in the Perimetric Area were detected at
 concentrations below background levels. Soil sample locations in the DDA and the
 Perimetric Area are indicated in Figure 4-1 from the July 2007 CSA.
- Groundwater samples and water table measurements have been collected from 12 monitoring wells screened at the water table, though some of these wells no longer exist due to RAM/IRA activity. Water level measurements indicate that DDA groundwater discharges to Cedar Swamp Brook. Of the nine remaining wells, two are upgradient of the DDA, two are in the center of the DDA, and five are downgradient of the DDA as indicated in Figure 4-2. Groundwater data confirmed the presence of metals, PAH, and EPH contamination within the boundary of the DDA. Metals and sporadic PAH/EPH were present at concentrations greater than background levels in groundwater samples collected from monitoring wells downgradient of the DDA. However, results of surface water and sediment sampling performed in Cedar Swamp Brook indicates that the impacted area does not extend to the brook.
- Soil sampling locations for asbestos were selected based on the delineated boundary of the DDA. Twenty samples were collected at or just beyond the DDA perimeter and 22 were collected from within the DDA area, as indicated in Figure 4-1. Separate samples were collected for surface (0-3 inches) and subsurface (3-24 inches) soils to evaluate the exposure potential associated with asbestos. Results confirmed the presence of asbestos at one location within the DDA boundary.

A-4 Number, Distribution, and Handling of Samples

- Sampling locations, as presented in the figures in this Phase II CSA Addendum and in the original July 2007 Phase II CSA, provide sufficient data for identifying the nature and extent of contaminants and conducting risk characterization. Discrete samples of soil, groundwater, surface water, and sediment were collected for laboratory analysis. Sample compositing was not performed prior to analysis.
- Soil sampling within the DDA was concentrated in areas where OHM was identified through test pits and soil borings. Soil sampling around the DDA was concentrated in the area of highest migration potential, along the steep slope between DDA and Cedar Swamp Brook. Five groundwater monitoring wells are positioned about equidistant along the downgradient (north) side of the Site, between DDA and Cedar Swamp Brook. Surface water and sediment were sampled at two upgradient and four downgradient locations.

A-5 Temporal Distribution of Samples

Groundwater conditions warrant monitoring over time since seasonal water table changes may affect movement of contaminants. The fill materials in the DDA have been in place for 30 or more years; therefore groundwater conditions are expected to be relatively stable. The RAM and IRA activities that concluded in 2006 removed source materials, and these excavations and placement of fill could alter localized groundwater movement. Groundwater monitoring continued for three additional sampling rounds in June/July and December 2007, and May 2008. These rounds -- coupled with the earlier sampling rounds in May 2005, June 2006, and August 2006 – provide sufficient data to



- assess temporal changes in groundwater conditions. Concentrations of contaminants appear to be stable or slowly diminishing over time, as indicated in Section 6.
- Soil concentrations are expected to diminish slowly over time based on migration of contaminants to groundwater, though this would be difficult to confirm since the same soil location cannot be sampled more than once. An evaluation of temporal changes in soil concentrations was not attempted. Surface water and sediment concentrations could change over time based on contaminant transport through groundwater; the temporal distribution of groundwater samples described above should be sufficient to predict changes in surface water or sediment.

A-6 Completeness of Data

No significant data gaps were found.

A-7 Inconsistency and Uncertainty

 No investigation results were identified that are inconsistent with the CSM or would suggest uncertainty for an RAO.

A-8 <u>Information Considered Unrepresentative</u>

- The earlier rounds of groundwater data were not included in the RC, as these are considered to be less representative of current conditions compared to the most recent data. The May 2008 groundwater sampling results were used in the RC as indicated in Appendix B. Concentrations of contaminants appear to be stable or slowly diminishing over time, as indicated in Section 6.
- Because an AUL prohibiting intrusive activities will be instituted, only soils from 0 to 3 feet below ground surface were included in the RC.
- Any soil samples that were removed during the RAM or IRA excavation activities were not included in the RC, since they are no longer representative of soil at the Site.

Based on the above analysis, the data collected in the DDA that were used to support the RC are considered to be representative of conditions in this exposure area. Data are representative both spatially and temporally.

11.2 Data Usability Assessment

B-1 List MCP Activities and Data Evaluated

- The MCP investigations that resulted in site characterization data are described in Section 4 of the July 2007 Phase II CSA and this Addendum. The investigations have included sampling soil, groundwater, surface water, and sediment for laboratory analysis; test pit visual observations; a geophysical survey to evaluate buried metal; and removal of drums and other wastes through RAM and IRA activities.
- The data that were evaluated for this DUA -- including sample locations, dates, and lab report numbers -- are presented in Appendix C of the July 2007 Phase II CSA and in Table 11-1 of this Addendum.

B-2 Appropriateness of Analytical Methods



 Only CAM-compliant data were used in the 2007 Phase II CSA and this Addendum. Appropriate analytical methods were selected to provide data that quantitatively supports an RAO. Method types are listed in the lab data packages provided in Appendix C of Weston's 2007 Phase II CSA, and in Appendix A of this Addendum.

B-3 Appropriateness of Reporting Limits

 Analytical methods with appropriate reporting limits were selected to provide data that quantitatively support an RAO. Method reporting limits are listed in the lab data packages provided in Appendix C of Weston's 2007 Phase II CSA, and in Appendix A of this Addendum.

B-4 Analytical Accuracy and Precision

- An indication of whether or not the laboratory met all CAM requirements and performance standards without qualification is presented in the MCP Certification Forms and Case Narratives at the beginning of each laboratory report.
- Any data qualifications that were necessary were indicated on the Baker Hughes Data Validation Qualifiers spreadsheet presented along with each laboratory report in Appendix C of the July 2007 Phase II CSA, or in Appendix D of this Addendum. Data qualifiers were also incorporated into the data tables in Sections 4 and 6 and the RC of the July 2007 CSA and this Addendum.

B-5 Field Data Usability

 Samples were collected in accordance with accepted environmental practice including the CAM Quality Assurance and Quality Control Guidelines. Matrix spike/matrix spike duplicates (MS/MSDs), equipment blanks, and field duplicates were collected and were employed for appropriate qualification of data.

B-6 Rejected Data

 Data rejected as a result of the evaluation process are qualified as "R" and will not be used to support an RAO. Acid extractable analytes were rejected in DD-MW-208-R01-001-X and DD-MW-208-R01-001-D due to low surrogate recovery as indicated in Appendix D.

Based on the above analysis, the data collected in the DDA are considered to be usable to support an RAO, subject to the specific qualifications for selected data. There are sufficient unqualified and usable data to support an RAO.



12.0 CONCLUSIONS [310 CMR 40.0835(4)(I)]

As described in Weston's 2007 Phase II CSA, a RAM, IRA, Phase II investigations, and a Method 1 RC had been conducted at the Site. A condition of NSR could not be documented for the Site, among other reasons due to the presence of ACM and asbestos-contaminated soils.

Subsequent sampling for asbestos in soil, and groundwater sampling, have further characterized the Site and are documented in this Phase II CSA Addendum. In addition, a Method 3 RC has been conducted for the DDA, which replaces the earlier conclusions of Weston's Method 1 RC. The Method 3 RC was required to evaluate potential exposure to asbestos in soil, because Method 1 standards for asbestos are not currently available. Also, a Method 3 RC was required to evaluate potential environmental risks. The following conclusions are made based on the original July 2007 Phase II CSA and this Addendum:

- Environmental conditions at the Site have been significantly improved as a result of RAM and IRA activities completed to date. Removal of the buried waste barrels and significant quantities of dioxin, asbestos, and metals-containing materials from the Site has resulted in reduction of the residual concentrations of these contaminants in soil at the Site.
- Imminent Hazards do not exist at this Site under present conditions and site controls.
- A condition of No Substantial Hazard to human health and the environment has been achieved at the Site.
- The hydrogeological investigations within the Site and vicinity indicates that the Cedar Swamp Brook adjacent to the DDA serves as the final discharge point for the groundwater in the local area; hence, it is concluded that contamination present in the soil and groundwater at the Site would not migrate across the Cedar Swamp Brook and impact the groundwater quality in off-site areas.
- Results from soil and groundwater samples at the Site indicate that the concentration of metals, dioxin TEQ, and EPH in soil at the Site are elevated compared to the background concentrations at the Site. Contamination by dioxin, metals, and EPH is primarily present in soil at 0 to 5 ft depth within the Western and Central Clearing Area.
- Asbestos is present in surface and subsurface soil at the Site.
- Surface water and sediment sampling in Cedar Swamp Brook indicate that potential contaminant migration pathways from surface runoff or groundwater discharge to surface water and/or sediment are not complete pathways.
- A condition of NSR in soil and groundwater has been achieved for current and reasonably foreseeable future exposures, assuming the implementation of an AUL prohibiting disruption of the ground surface.
- A condition of NSR has been achieved for the environment as well.
- Background conditions have not been met for PAHs and for the metals barium, cadmium, chromium, mercury, nickel, vanadium, and zinc in Site groundwater.
- Background conditions exist in soil based on sampling performed in the Perimetric Area between the Site and Cedar Swamp Brook.

Based on the results of this Phase II CSA, a condition of NSR has been achieved at the Site.

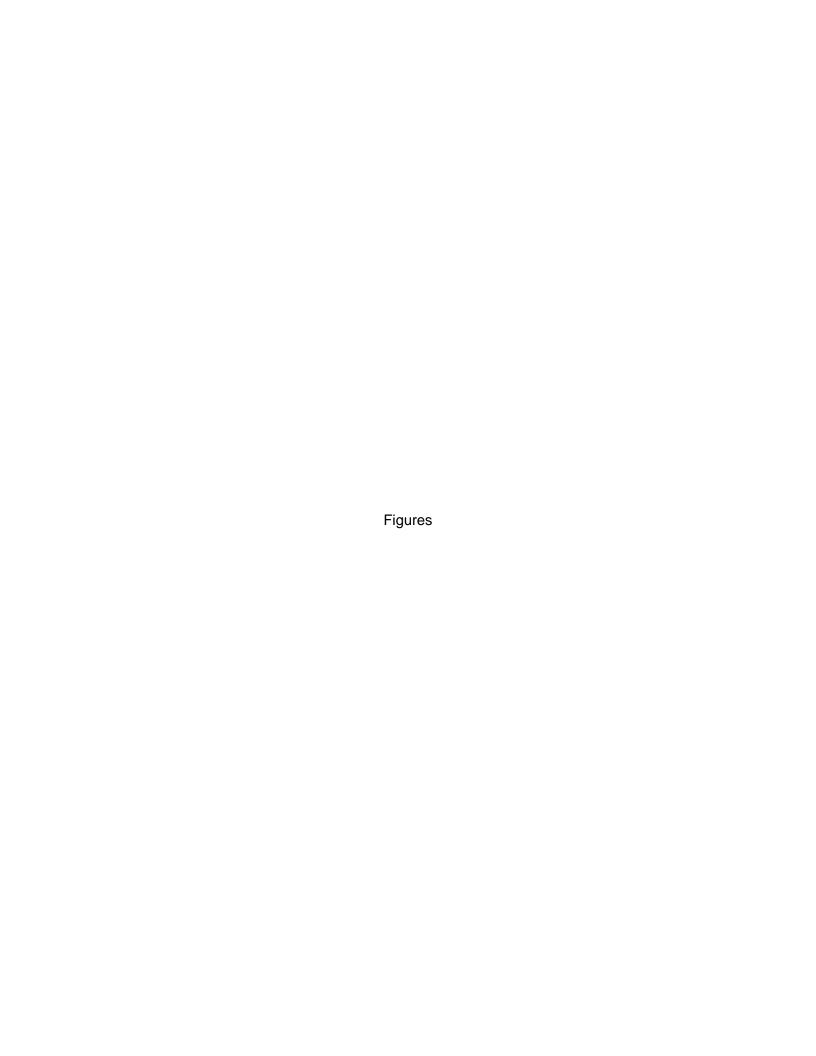


13.0 REFERENCES

MADEP, 2003. Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data. Bureau of Waste Site Cleanup, Rev 3.1, May 22.

U.S. EPA, 1996. Method 1669 Sampling Ambient Water for Trace Metals at EPA Water Quality Criteria Levels. July

Weston, 2007. Phase II Comprehensive Site Assessment Report, Demolition Debris Area, RTN 3-3023105. 30 July



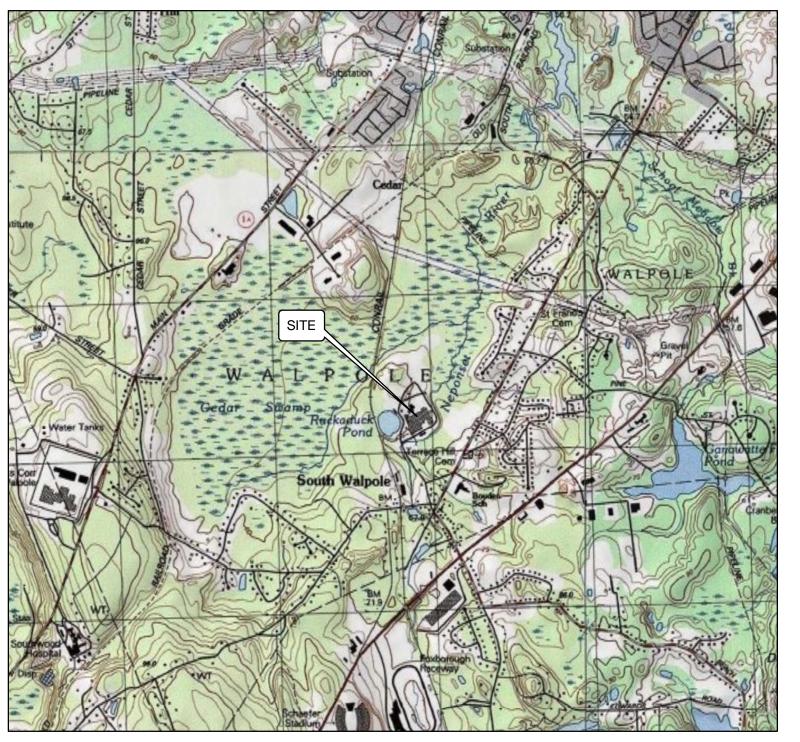


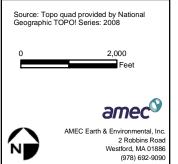
FIGURE 1-1

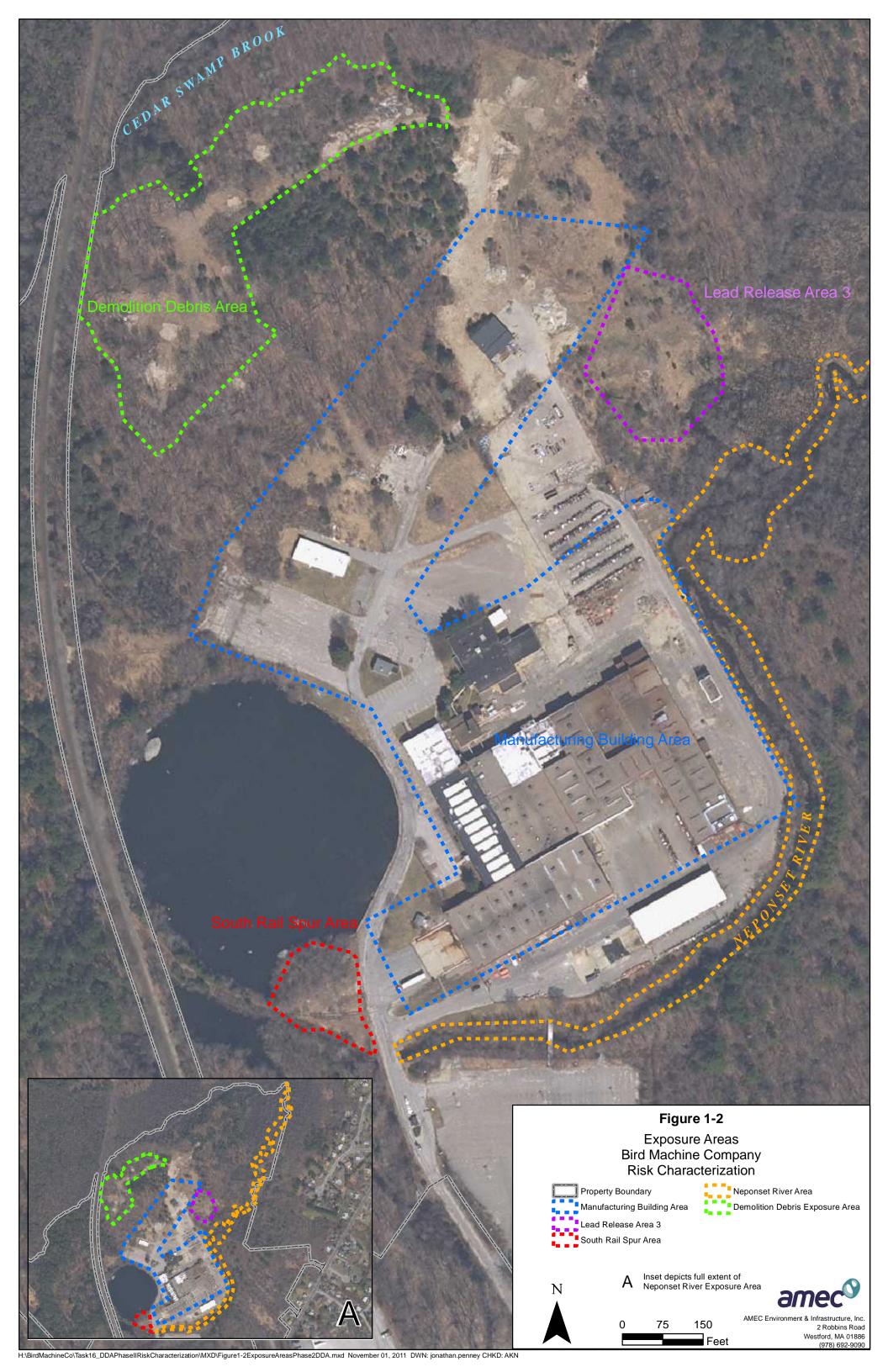
SITE LOCATION MAP

Bird Machine Inc. Company

100 Neponset Street Walpole, MA







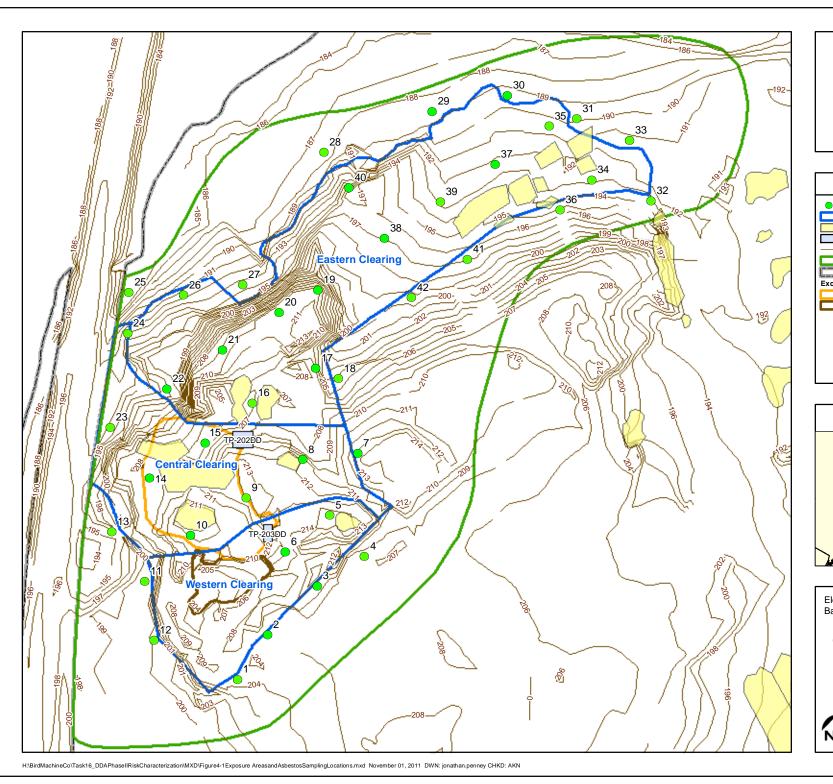


FIGURE 4-1

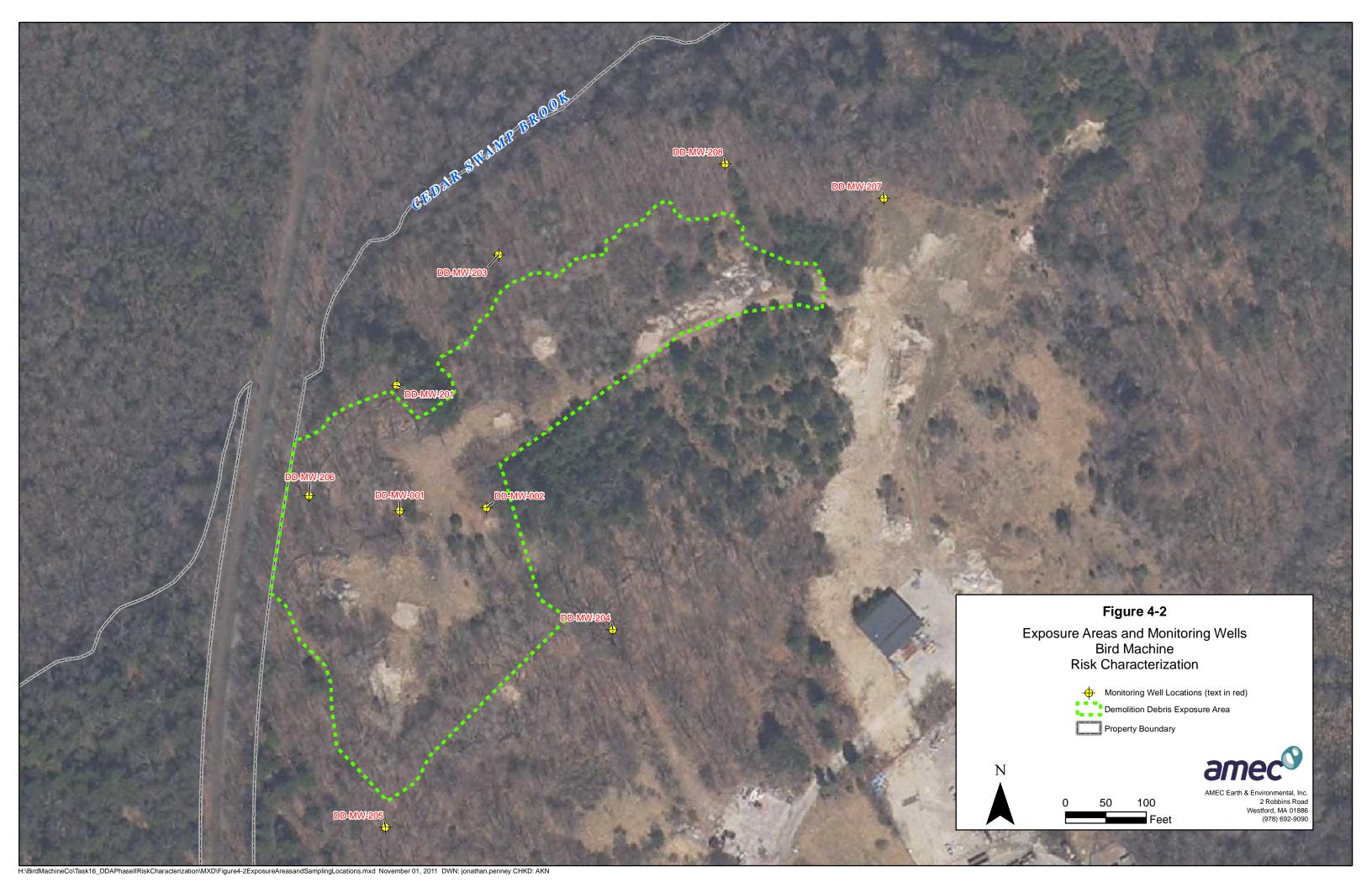
Asbestos In Soil Sampling Locations

Bird Machine Co. Walpole, MA









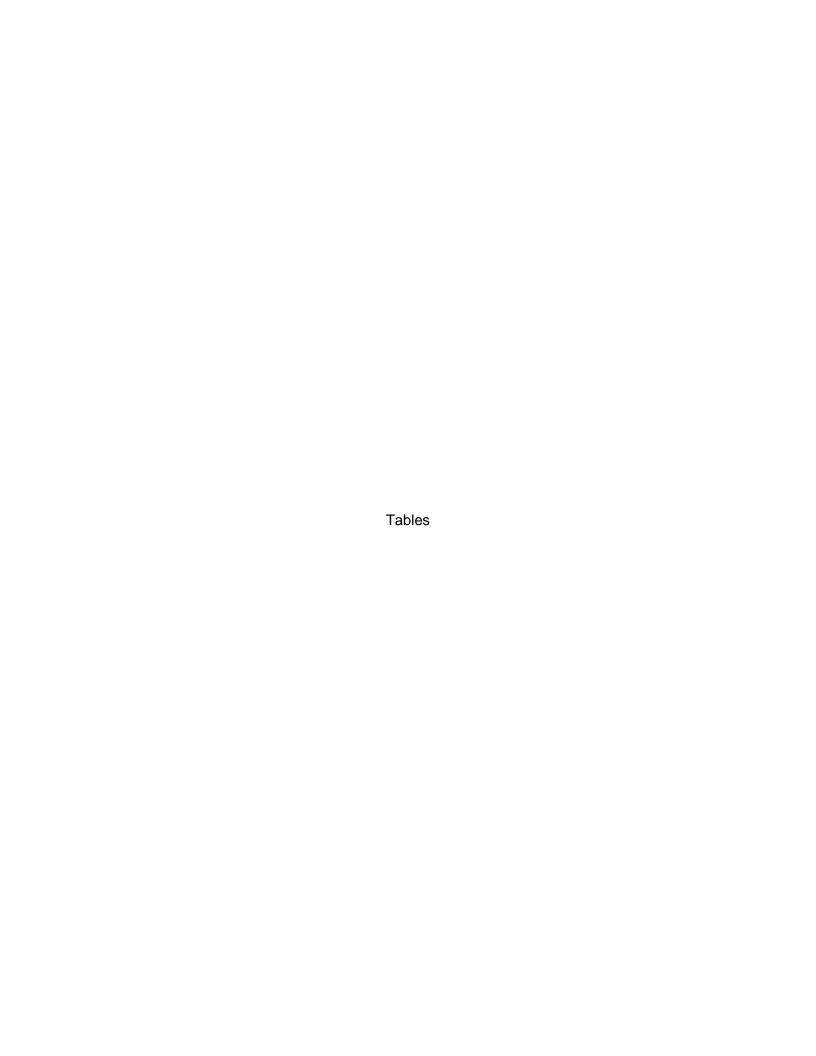


Table 4-1 Asbestos Sampling and Analysis Plan Phase II CSA Addendum Demolition Debris Area Bird Machine Company - Walpole, MA

Sample ID (1)	Fill Observed (2)	PLM (3)	TEM ⁽⁴⁾
1 A & B	No	X	
1A-DUP	No	X	
2 A & B	Yes	X	
2A-DUP	Yes	X	
3 A & B	Yes	X	
3A-DUP	Yes	Х	
4 A & B	No	X	
4A-DUP	No	X	
5 A & B	No	X	
5A-DUP	No	X	
6 A & B	Yes	X	
6A-DUP	Yes	X	
7 A & B	Yes	X	
7A-DUP	Yes	X	
8AS	Yes	X	
8AS - TEM	Yes	X	
8 A & B	Yes	X	
8 A & B - TEM	Yes	X	X
8A - TEM DUP	Yes	X	X
9 A & B	No	X	
9A-DUP	No	X	
10 A & B	Yes	X	
11 A & B	No	X	
12 A & B	No	X	
13 A & B	No	X	
14 A & B	Yes	X	
14 A & B - TEM	Yes	X	Χ
15 A & B	Yes	X	Α
16 A & B	No	X	
17 A & B	Yes	X	
18 A & B	No Var	X	
19 A & B	Yes	X	
20 A & B	Yes	X	
21 A & B	Yes	X	
22 A & B	Yes	X	
23 A & B	No	X	
24 A & B	No	X	
25 A & B	No	X	
26 A & B	No	X	
27 A & B	No	X	
28 A & B	No	Х	
29 A & B	Yes	X	
30 A & B	Yes	X	
31 A & B	No	X	
32 A & B	No	X	
33 A & B	No	X	
33 A & B	No	X	X
34 A & B	No	X	
35 A & B	No	X	
36 A & B	No	X	
		X	
36 A & B	No You		
37 A & B	Yes	X	
38 A & B	Yes	X	
39 A & B	No	X	
40 A & B	Yes	X	
41 A & B	No	X	
42 A & B	No	X	

Notes:

(1)Samples were collected from two depths at each location, with the exception of location 8. For all samples, "A" designates samples collected from the 0-3" interval while "B" designates samples collected from 3-24" interval. A surface soil sample was also collected at location 8 which is designated by "AS" (2) Visual characterization of samples as from fill material or native material

- (3) "X" indicates sample was analyzed for total asbestos via Polarized Light Microscopy (PLM)
- (4) "X" indicates sample was analyzed for asbestos via Transmission Electron Microscopy (TEM)

Table 4-2 **Asbestos Soil Sampling Results** Phase II CSA Addendum **Demolition Debris Area** Bird Machine Company - Walpole, MA

Sample ID	Depth	PLM Result (1)		TEM Result (2)	
8AS -AR	Surface	1.9	Chrysotile		
8AS-TEM	Surface	1.7 5.2	Chrysotile Crocidolite		
8A	0-3"	1.3	Chrysotile		
8A-TEM	0-3"	trace	Chrysotile	ND	
8A-TEM-DUP	0-3"	trace	Chrysotile	trace	Chrysotile
8B-TEM	3-24"	trace	Chrysotile	trace	Chrysotile

Notes:

- (1) Results for samples analysed for total asbestos via Polaized Light Microscopy (PLM)(2) Results for samples analysed for total asbestos via Transmission Electron Microscopy (TEM)
- All samples were run for PLM analysis, while a smaller subset was analyzed via TEM
- ND Sample was non-detect for asbestos

Table 4-3 Results of Asbestos Elutriator Analysis Phase II CSA Addendum Demolition Debris Area Bird Machine Company - Walpole, MA

Sample ID	Depth	Elutriator Result Total Asbestos Fibers (MFG)
8AS -AR	Surface	15.1
8AS-TEM	Surface	82.9
8A-TEM	0-3"	11.7

Notes:

MFG - million fibers per gram Results for samples analyzed for total asbestos via Elutriator Method Analysis

		ation:	DD-MW-001		DD-MW-002		DD-MW-20		DD-MW-203		DD-MW-204		DD-MW-205		DD-MW-206		DD-MW-207	
	Field Samp Sample		DD-MW-001-R0 6/5/2007	2-X	DD-MW-002-R0 6/5/2007)1-X	DD-MW-201-R0 6/5/2007	04-X	DD-MW-203-R0 6/5/2007	4-X	DD-MW-204-R0 6/5/2007)2-X	DD-MW-205-R0 6/6/2007	3-X	DD-MW-206-R0 6/5/2007	4-X	DD-MW-207-R0 6/6/2007)4-X
CAS Number		epth: Units	11.6 - 16.6 fee Result	et Q	9.8 - 14.8 fee Result	et Q	4 - 14 feet Result	Ю	4.5 - 12.5 fee Result	t Q	7.5 - 13.5 fee Result	et Q	5 - 11 feet Result	Q	5 - 11 feet Result	Q	5 - 15 feet Result	Q
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	5.1	U	5.1	U	5.3	U	5.2	U	5.1	U	5.1	U	5.2	U	5.1	U
95-50-1 541-73-1	1,2-DICHLOROBENZENE 1,3-DICHLOROBENZENE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
106-46-7 95-95-4	1,4-DICHLOROBENZENE 2,4,5-TRICHLOROPHENOL	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	UJ	5.1 5.1	U	5.1 5.1	U UJ	5.2 5.2	UJ	5.1 5.1	U UJ
88-06-2 120-83-2	2,4,6-TRICHLOROPHENOL 2.4-DICHLOROPHENOL	ug/l ug/l	5.1 5.1	UJ	5.1 5.1	UJ	5.3 5.3	UJ	5.2 5.2	UJ	5.1 5.1	UJ	5.1 5.1	UJ	5.2 5.2	UJ	5.1 5.1	UJ
105-67-9 51-28-5	2,4-DIMETHYLPHENOL 2.4-DINITROPHENOL	ug/l ug/l	5.1 5.1	UJ	5.1 5.1	UJ	5.3 5.3	UJ	5.2	UJ	5.1 5.1	UJ	5.1 5.1	UJ	5.2	UJ	5.1 5.1	UJ
121-14-2	2,4-DINITROTOLUENE	ug/l	5.1	U	5.1	U	5.3	U	5.2	U	5.1	U	5.1	U	5.2	U	5.1	U
606-20-2 91-58-7	2,6-DINITROTOLUENE 2-CHLORONAPHTHALENE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
95-57-8 91-57-6	2-CHLOROPHENOL 2-METHYLNAPHTHALENE	ug/l ug/l	5.1 1	UJ	5.1 1	UJ	5.3 1.1	UJ	5.2 1	UJ	5.1 1	UJ	5.1 1	UJ	5.2 1	UJ	5.1 1	UJ
95-48-2 95-48-7	2-METHYLPHENOL 2-METHYLPHENOL (O-CRESOL)	ug/l ug/l	5.1 NA	UJ	5.1 NA	UJ	5.3 NA	UJ	5.2 NA	UJ	5.1 NA	UJ	5.1 NA	UJ	5.2 NA	UJ	5.1 NA	UJ
88-75-5 106-44-5	2-NITROPHENOL 3&4-METHYLPHENOL	ug/l ug/l	5.1 5.1	UJ	5.1 5.1	UJ	5.3 5.3	UJ	5.2 5.2	UJ	5.1 5.1	UJ	5.1 5.1	UJ	5.2 5.2	UJ	5.1 5.1	UJ
91-94-1	3,3-DICHLOROBENZIDINE	ug/l	5.1	U	5.1	U	5.3	U	5.2	U	5.1	U	5.1	U	5.2	U	5.1	U
101-55-3 106-47-8	4-BROMOPHENYL PHENYL ETHER 4-CHLOROANILINE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
100-02-7 83-32-9	4-NITROPHENOL ACENAPHTHENE	ug/l ug/l	5.1 1	UJ	5.1 1	UJ	5.3 1.1	UJ	5.2 1	UJ	5.1 1	UJ	5.1 1	UJ	5.2 1	UJ	5.1 1	UJ
208-96-8 98-86-2	ACENAPHTHYLENE ACETOPHENONE	ug/l ug/l	0.3 5.1	U	0.3 5.1	U	0.32 5.3	U	0.31 5.2	U	0.31 5.1	U	0.3 5.1	U	0.31 5.2	U	0.3 5.1	U
62-53-3	ANILINE	ug/l	51	UJ	51	UJ	53	UJ	52	UJ	51	UJ	51	UJ	52	UJ	51	UJ
120-12-7 103-33-3	ANTHRACENE AZOBENZENE	ug/l ug/l	1 5.1	U	5.1	U	1.1 5.3	U	1 5.2	U	1 5.1	U	1 5.1	U	1 5.2	U	1 5.1	U
56-55-3 50-32-8	BENZO(A)ANTHRACENE BENZO(A)PYRENE	ug/l ug/l	0.3 0.2	U	0.3 0.2	U	0.32 0.21	U	0.31 0.21	U	0.31 0.2	U	0.3 0.2	U	0.31 0.21	U	0.22 0.2	J
205-99-2 191-24-2	BENZO(B)FLUORANTHENE BENZO(G,H,I)PERYLENE	ug/l	0.3 0.51	U	0.3 0.51	U	0.32 0.53	U	0.31 0.52	U	0.31 0.51	U	0.3	U	0.31 0.52	U	0.3 0.51	U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3	U	0.3	Ü	0.32	U	0.31	U	0.31	Ū	0.3	U	0.31	U	0.3	U
111-91-1 111-44-4	BIS(2-CHLOROETHOXY)METHANE BIS(2-CHLOROETHYL) ETHER	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
108-60-1 117-81-7	BIS(2-CHLOROISOPROPYL)ETHER BIS(2-ETHYLHEXYL) PHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
85-68-7 218-01-9	BUTYLBENZYLPHTHALATE CHRYSENE	ug/l ug/l	5.1 1	U	5.1 1	U	5.3 1.1	U	5.2 1	U	5.1	U	5.1 1	U	5.2 1	U	5.1 0.24	U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.51	U	0.51	Ü	0.53	Ü	0.52	U	0.51	U	0.51	U	0.52	U	0.51	U
132-64-9 84-66-2	DIBENZOFURAN DIETHYL PHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
131-11-3 84-74-2	DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.3 5.3	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U
117-84-0 206-44-0	DI-N-OCTYL PHTHALATE FLUORANTHENE	ug/l ug/l	5.1 1	U	5.1 1	U	5.3 1.1	U	5.2 1	U	5.1 1	U	5.1 1	U	5.2 1	U	0.39 0.28	J
86-73-7	FLUORENE	ug/l	1	U	1	Ü	1.1	U	1	U	1	Ū	1	U	1	U	1	Ü
118-74-1 87-68-3	HEXACHLOROBENZENE HEXACHLOROBUTADIENE	ug/l ug/l	1 0.61	U	0.61	U	1.1 0.63	U	1 0.62	U	1 0.61	U	1 0.61	U	1 0.63	U	1 0.61	U
67-72-1 193-39-5	HEXACHLOROETHANE INDENO(1,2,3-CD)PYRENE	ug/l ug/l	5.1 0.51	U	5.1 0.51	U	5.3 0.53	U	5.2 0.52	U	5.1 0.51	U	5.1 0.51	U	5.2 0.52	U	5.1 0.51	U
78-59-1 91-20-3	ISOPHORONE NAPHTHALENE	ug/l ug/l	5.1 1	U	5.1 1	U	5.3 1.1	U	5.2 1	U	5.1 1	U	5.1 1	U	5.2 1	U	5.1 1	U
98-95-3 106-44-5	NITROBENZENE P-CRESOL	ug/l	5.1 NA	Ü	5.1 NA	Ü	5.3 NA	U	5.2 NA	Ü	5.1 NA	Ü	5.1 NA	Ü	5.2 NA	Ü	5.1 NA	Ü
87-86-5	PENTACHLOROPHENOL	ug/l ug/l	5.1	UJ	5.1	UJ	5.3	UJ	5.2	UJ	5.1	UJ	5.1	UJ	5.2	UJ	5.1	UJ
85-01-8 108-95-2	PHENANTHRENE PHENOL	ug/l ug/l	0.2 5.1	U	0.2 5.1	U UJ	0.21 5.3	UJ	0.21 5.2	UJ	0.2 5.1	U	0.2 5.1	U UJ	0.21 5.2	UJ	0.2 5.1	U
129-00-0	PYRENE	ug/l	5.1	U	5.1	U	5.3	U	5.2	U	5.1	U	5.1	U	5.2	U	0.29	J
35822-46-9 67562-39-4	1,2,3,4,6,7,8-HpCDD 1,2,3,4,6,7,8-HpCDF	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
55673-89-7 39227-28-6	1,2,3,4,7,8,9-HpCDF 1,2,3,4,7,8-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\perp
70648-26-9 57653-85-7	1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\perp
57117-44-9 19408-74-3	1,2,3,6,7,8-HxCDF 1,2,3,7,8,9-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\perp
72918-21-9 40321-76-4	1,2,3,7,8,9-HxCDF 1,2,3,7,8-PeCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
57117-41-6 60851-34-5	1,2,3,7,8-PeCDF 2,3,4,6,7,8-HxCDF	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\vdash
57117-31-4 1746-01-6	2,3,4,7,8-PeCDF 2,3,7,8-TCDD	pg/l pg/l	NA NA	E	NA NA	E	NA NA	E	NA NA	H	NA NA	H	NA NA	Ħ	NA NA	H	NA NA	且
51207-31-9 37871-00-4	2,3,7,8-TCDF HPCDD (TOTAL)	pg/l pg/l	NA NA	E	NA NA	E	NA NA	F	NA NA	H	NA NA	Ħ	NA NA	H	NA NA	H	NA NA	彐
38998-75-3 34465-46-8	HPCDF (TOTAL) HXCDD (TOTAL)	pg/l pg/l	NA NA		NA NA		NA NA	F	NA NA	H	NA NA	Н	NA NA	H	NA NA		NA NA	F
55684-94-1 3268-87-9	HXCDF (TOTAL) OCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA	H	NA NA		NA NA	H	NA NA		NA NA	H
39001-02-0 36088-22-9	OCDF PECDD (TOTAL)	pg/l pg/l	NA NA		NA NA		NA NA		NA NA	H	NA NA	H	NA NA		NA NA		NA NA	Ħ
30402-15-4 41903-57-5	PECDF (TOTAL) TCDD (TOTAL)	pg/l pg/l	NA NA		NA NA		NA NA		NA NA	H	NA NA	H	NA NA		NA NA		NA NA	Ħ
55722-27-5	TCDF (TOTAL)	pg/l	NA		NA		NA		NA		NA		NA		NA		NA	
91-57-6 83-32-9	2-METHYLNAPHTHALENE ACENAPHTHENE	ug/l ug/l	1 1	U	1.1 1.1	U	1.1 1.1	U	1 1	U U	1 1	U	1 1	U	1 1	U	1.1 1.1	U
208-96-8 120-12-7	ACENAPHTHYLENE ANTHRACENE	ug/l ug/l	0.3 1	U	0.32 1.1	U	0.32 1.1	U	0.3 1	U	0.3 1	U	0.3 1	U	0.31 1	U	0.32 1.1	U
56-55-3 50-32-8	BENZO(A)ANTHRACENE BENZO(A)PYRENE	ug/l ug/l	0.3 0.2	U	0.32 0.21	U	0.32 0.21	U	0.3 0.2	U	0.3 0.2	U	0.3 0.2	U	0.31 0.21	U	0.32 0.21	U
205-99-2 191-24-2	BENZO(B)FLUORANTHENE BENZO(G,H,I)PERYLENE	ug/l ug/l	0.3 0.4	U	0.32 0.42	U	0.32 0.42	U	0.3 0.4	U	0.3 0.4	U	0.3 0.4	U	0.31 0.41	U	0.32 0.42	U
207-08-9 EPH1122	BENZO(K)FLUORANTHENE C11-C22 AROMATICS, ADJUSTED	ug/l ug/l	0.3 100	U	0.32 110	U	0.32 110	U	0.3 100	U	0.3 100	U	0.3 100	U	0.31 100	U	0.32 110	U
EPH1122 EPH1122	C11-C22 AROMATICS, ADJUSTED C11-C22 AROMATICS, UNADJUSTED	ug/l ug/l	NA 100	U	NA 110	U	NA 110	U	NA 100	U	NA 100	U	NA 100	U	NA 100	U	NA 110	U
EPH1936 EPH1936	C19-C36 ALIPHATICS C19-C36 ALIPHATICS (EPH-LL)	ug/l ug/l	100 NA	Ü	110 110 NA	Ü	110 110 NA	Ü	100 NA	Ü	100 100 NA	Ü	100 NA	Ü	100 100 NA	Ü	110 NA	Ü
VPH912 VPH912	C9-C18 ALIPHATICS (EPH-LL) C9-C18 ALIPHATICS (EPH-LL)	ug/l ug/l	100 NA	U	110 NA	U	110 NA	U	100 NA	U	100 NA	U	100 NA	U	100 NA	U	110 NA	U
218-01-9	CHRYSENE DIBENZ(A,H)ANTHRACENE	ug/l ug/l	1 0.4	U	1.1 0.42	U	1.1 0.42	U	1 0.4	U	1 0.4	U	1 0.4	U	1 0.41	U	1.1 0.42	U
53-70-3	IDIDEIXE(A,II)AN IIINACENE	uy/I			0.42				∪.⊶	J	U. 4	U		J	U. + I		U.4Z	
53-70-3 206-44-0 86-73-7	FLUORANTHENE	ug/l	1	U	1.1	U	1.1	U	1	U	1	U	1	U	1	U	1.1	U
206-44-0 86-73-7 193-39-5	FLUORENE INDENO(1,2,3-CD)PYRENE	ug/l ug/l	1 0.4	U	1.1 0.42	U	1.1 0.42	U	1 0.4	U	1 0.4	U	1 0.4	U	1 0.41	U	1.1 0.42	U
206-44-0 86-73-7	FLUORENE	ug/l	1	U	1.1	Ū	1.1	U	1	Ü	1	Ū	1	U	1	Ü	1.1	U

	Field Samp		DD-MW-001-R0		DD-MW-002-R		DD-MW-201-R0)4-X	DD-MW-203-R0		DD-MW-204 DD-MW-204-R0		DD-MW-205 DD-MW-205-R0		DD-MW-206 DD-MW-206-R0		DD-MW-207-R	
CAS Number		Date: Depth: Units	6/5/2007 11.6 - 16.6 fe Result	et Q	6/5/2007 9.8 - 14.8 fee Result	et Q	6/5/2007 4 - 14 feet Result	Q	6/5/2007 4.5 - 12.5 fee Result	t Q	6/5/2007 7.5 - 13.5 fee Result	t Q	6/6/2007 5 - 11 feet Result	Q	6/5/2007 5 - 11 feet Result	Q	6/6/2007 5 - 15 feet Result	t Q
7440-36-0	ANTIMONY ANTIMONY (DISSOLVED)	ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-36-0 7440-38-2 7440-38-2	ARSENIC ARSENIC (DISSOLVED)	ug/l ug/l ug/l	NA 2	U	NA 2	U	NA 0.81	ı	NA 1	U	NA 1	U	NA 2	U	NA 1	U	NA 5.7	V
7440-39-3 7440-39-3	BARIUM BARIUM (DISSOLVED)	ug/l ug/l	NA 47	V	NA 32	V	NA 44	V	NA 12	V	NA 17	V	NA 18	V	NA 26	V	NA 38	V
7440-41-7 7440-41-7	BERYLLIUM BERYLLIUM (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-43-9 7440-43-9	CADMIUM CADMIUM (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-47-3 7440-47-3	CHROMIUM CHROMIUM (DISSOLVED)	ug/l ug/l	NA 35	V	NA 7.3	V	NA 25	V	NA 16	V	NA 11	V	NA 8.4	V	NA 23	V	NA 4.9	V
7440-50-8 7440-50-8	COPPER (DISSOLVED)	ug/l ug/l	NA NA	Ľ	NA NA	Ļ	NA NA	Ľ	NA NA	Ľ	NA NA		NA NA		NA NA	Ė	NA NA	Ť
7439-92-1 7439-92-1	LEAD LEAD (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7439-97-6 7439-97-6	MERCURY MERCURY (DISSOLVED)	ug/l ug/l	NA 0.2	U	NA 0,2	U	NA 0.2	U	NA 0.2	U	NA 0.2	U	NA 0.2	IJ	NA 0,2	IJ	NA 0.2	U
7440-02-0 7440-02-0	NICKEL NICKEL (DISSOLVED)	ug/l ug/l	NA 7.2	V	NA 2.5	V	NA 1.3	V	NA 1.1	V	NA 1.1	V	NA 1.4	V	NA 0.91	J.	NA 1	U
7782-49-2 7782-49-2	SELENIUM SELENIUM (DISSOLVED)	ug/l ug/l	NA 2	U	NA 2	U	NA 9.3	J	NA 1	U	NA 1	U	NA 2	U	NA 1	U	NA 1	U
7440-22-4 7440-22-4	SILVER SILVER (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-28-0 7440-28-0	THALLIUM THALLIUM (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-62-2 7440-62-2	VANADIUM VANADIUM (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-66-6 7440-66-6	ZINC ZINC (DISSOLVED)	ug/l ug/l	NA 5	U	NA 5	U	NA 2.5	U	NA 2.4	J	NA 2.1	J	NA 5	U	NA 2.7	٧	NA 2.5	U
630-20-6	1,1,1,2-TETRACHLOROETHANE	ug/l	NA	Ě	NA	Ĕ	NA	Ĕ	NA		NA	Ě	NA		NA	Ė	NA	Ť
71-55-6 79-34-5	1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE	ug/l ug/l	NA NA	Ė	NA NA	Ē	NA NA	Ē	NA NA	E	NA NA	E	NA NA	Ы	NA NA	H	NA NA	士
79-00-5 75-34-3	1,1,2-TRICHLOROETHANE 1,1-DICHLOROETHANE	ug/l ug/l	NA NA	Ē	NA NA	Ē	NA NA	Ē	NA NA	E	NA NA	E	NA NA	H	NA NA		NA NA	上
75-35-4 563-58-6	1,1-DICHLOROETHENE 1,1-DICHLOROPROPENE	ug/l ug/l	NA NA	Ē	NA NA	Ē	NA NA	Ē	NA NA	Ē	NA NA	Ē	NA NA	Ы	NA NA	H	NA NA	\pm
87-61-6 96-18-4	1,2,3-TRICHLOROBENZENE 1,2,3-TRICHLOROPROPANE	ug/l ug/l	NA NA	E	NA NA	Ē	NA NA	Ē	NA NA	Ē	NA NA	E	NA NA		NA NA	H	NA NA	士
120-82-1 95-63-6	1,2,4-TRICHLOROBENZENE 1,2,4-TRIMETHYLBENZENE	ug/l ug/l	NA NA	Ė	NA NA	Ē	NA NA	Ē	NA NA	E	NA NA	E	NA NA	Ы	NA NA	H	NA NA	士
96-12-8 95-50-1	1,2-DIBROMO-3-CHLOROPROPANE 1,2-DICHLOROBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
107-06-2 78-87-5	1,2-DICHLOROETHANE 1,2-DICHLOROPROPANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
108678 541-73-1	1,3,5-TRIMETHYLBENZENE 1,3-DICHLOROBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
142-28-9 106-46-7	1,3-DICHLOROPROPANE 1,4-DICHLOROBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
123-91-1 594-20-7	1,4-DIOXANE 2,2-DICHLOROPROPANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
78-93-3 95-49-8	2-BUTANONE 2-CHLOROTOLUENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
591-78-6 106-43-4	2-HEXANONE 4-CHLOROTOLUENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	lacksquare
99-87-6 108-10-1	4-ISOPROPYLTOLUENE 4-METHYL-2-PENTANONE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
67-64-1 71-43-2	ACETONE BENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	士
108-86-1 75-27-4	BROMOBENZENE BROMODICHLOROMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	士
75-25-2 74-83-9	BROMOFORM BROMOMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	士
75-15-0 56-23-5	CARBON DISULFIDE CARBON TETRACHLORIDE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
108-90-7 74-97-5	CHLOROBENZENE CHLOROBROMOMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	士
75-00-3 67-66-3	CHLOROETHANE CHLOROFORM	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
74-87-3 156-59-2	CHLOROMETHANE CIS-1,2-DICHLOROETHENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
10061-01-5 124-48-1	CIS-1,3-DICHLOROPROPENE DIBROMOCHLOROMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	士
74-95-3 75-71-8	DIBROMOMETHANE DICHLORODIFLUOROMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
60-29-7 108-20-3	DISOPROPYL ETHER (DIPE)	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA		NA NA	Н	NA NA	Ы	NA NA	士
100-41-4 106-93-4	ETHYLBENZENE ETHYLENE DIBROMIDE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	H	NA NA		NA NA	丰
637-92-3 87-68-3	HEXACHLOROBUTADIENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	H	NA NA	H	NA NA	丰
98-82-8	ISOPROPYL ETHER ISOPROPYLBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA		NA NA	H	NA NA		NA NA	#
75-09-2	M,P-XYLENES METHYLENE CHLORIDE METHYL T BUTYL ETHER (MTRE)	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	H	NA NA		NA NA		NA NA		NA NA	#
1634-04-4 91-20-3	METHYL-T-BUTYL ETHER (MTBE) NAPHTHALENE	ug/l ug/l	NA NA	F	NA NA	1	NA NA	F	NA NA	Ħ	NA NA		NA NA	H	NA NA		NA NA	#
104-51-8 103-65-1	N-BUTYLBENZENE N-PROPYLBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA		NA NA	H	NA NA		NA NA	丰
95-47-6 135-98-8	O-XYLENE SEC-BUTYLBENZENE	ug/l ug/l	NA NA		NA NA	1	NA NA		NA NA	H	NA NA		NA NA	H	NA NA		NA NA	#
100-42-5 994-05-8	STYRENE TERT-AMYL METHYL ETHER (TAME)	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA		NA NA	H	NA NA		NA NA	丰
98-06-6 127-18-4	TERT-BUTYLBENZENE TETRACHLOROETHYLENE	ug/l ug/l	NA NA	H	NA NA	-	NA NA NA		NA NA	Ħ	NA NA		NA NA	H	NA NA NA		NA NA	#
109-99-9 108-88-3	TETRAHYDROFURAN TOLUENE TRANS 1.2 DICHLOROETHENE	ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA		NA NA NA	H	NA NA NA		NA NA NA	#
156-60-5 10061-02-6 79-01-6	TRANS-1,2-DICHLOROETHENE TRANS-1,3-DICHLOROPROPENE TRICHLOROETHYLENE (TCE)	ug/l ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA	H	NA NA NA	#
75-69-4	TRICHLOROFLUOROMETHANE	ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA		NA NA NA	H	NA NA NA		NA NA NA	#
75-01-4 71-43-2	VINYL CHLORIDE BENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丰
VPH58 VPH58	C5-C8 ALIPHATICS, ADJUSTED C5-C8 ALIPHATICS. UNADJUSTED	ug/l ug/l	NA NA NA	H	NA NA NA	1	NA NA NA		NA NA NA	Ħ	NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	丰
VPH910 VPH912	C9-C10 AROMATICS C9-C12 ALIPHATICS, ADJUSTED	ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	#
VPH912 100-41-4	C9-C12 ALIPHATICS, ADJUSTED C9-C12 ALIPHATICS, UNADJUSTED ETHYLBENZENE	ug/l ug/l	NA NA NA	H	NA NA NA	1	NA NA NA		NA NA NA	Ħ	NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	丰
1634-04-4	M,P-XYLENES METHYL-T-BUTYL ETHER (MTBE)	ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	#
91-20-3	NAPHTHALENE O-XYLENE	ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA	H	NA NA NA	#
95-47-6	IO-VIETNE	ug/I	INA	4	NA NA	1	INA	1	i INA	1	INA	1	INA		INA	i	INA	

Notes:
ug/L - micrograms per liter
V - Valid result, no qualification needed
U - not detected, laboratory reporting limit listed
J - concentration (or reporting limit) is estimated
DDA - Demolition Debris Area
NA - Not analyzed

	Sta	ation:	DD-MW-208	3	DD-MW-20	8	DD-MW-208	3	DD-MW-201		DD-MW-203		DD-MW-207		DD-MW-207	,	DD-MW-20	18
	Field Samp Sample		DD-MW-208-R01- 6/25/2007	001-	DD-MW-208-R01 6/25/2007	-001-	DD-MW-208-R0 7/23/2007	02-X	DD-MW-201-R0 12/11/2007	15-X	DD-MW-203-R09	5-X	DD-MW-207-R0 12/11/2007	15-X	DD-MW-207-R0 12/11/2007)5-D	DD-MW-208-R 12/11/2007	
CAS Number		epth: Units	3 - 13 feet Result	Q	3 - 13 feet Result	Q	3 - 13 feet Result	Q	4 - 14 feet Result	Q	4.5 - 12.5 feet Result	t Q	5 - 15 feet Result	Q	5 - 15 feet Result	Q	3 - 13 feet Result	
120-82-1 95-50-1	1,2,4-TRICHLOROBENZENE 1,2-DICHLOROBENZENE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
541-73-1	1,3-DICHLOROBENZENE	ug/l	5.1	U	5.1	U	5.2	U	5.1	U	5.1	U	5.1	U	5.5	U	5.1	U
106-46-7 95-95-4	1,4-DICHLOROBENZENE 2,4,5-TRICHLOROPHENOL	ug/l ug/l	5.1 5.1	U R	5.1 5.1	U R	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
88-06-2 120-83-2	2,4,6-TRICHLOROPHENOL 2,4-DICHLOROPHENOL	ug/l ug/l	5.1 5.1	R R	5.1 5.1	R R	5.2 5.2	UJ	5.1 5.1	\Box	5.1 5.1	U	5.1 5.1	\subset	5.5 5.5	\Box	5.1 5.1	U
105-67-9	2,4-DINITROPHENOL	ug/l	5.1	R	5.1	R	5.2 5.2	UJ	5.1	UJ	5.1	U	5.1	U	5.5 5.5	U	5.1	U
51-28-5 121-14-2	2,4-DINITROTOLUENE	ug/l ug/l	5.1 5.1	R U	5.1 5.1	R U	5.2	U	5.1 5.1	U	5.1 5.1	UJ	5.1 5.1	UJ	5.5	U	5.1 5.1	U
91-58-7	2,6-DINITROTOLUENE 2-CHLORONAPHTHALENE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
95-57-8 91-57-6	2-CHLOROPHENOL 2-METHYLNAPHTHALENE	ug/l ug/l	5.1 1	R	5.1 1	R	5.2 1	UJ	5.1 1	U	5.1 1	U	5.1 1	U	5.5 1.1	U	5.1 1	U
95-48-2	2-METHYLPHENOL	ug/l	5.1	R	5.1	R	5.2	UJ	5.1	U	5.1	Ü	5.1	U	5.5	U	5.1	U
95-48-7 88-75-5	2-METHYLPHENOL (O-CRESOL) 2-NITROPHENOL	ug/l ug/l	NA 5.1	R	NA 5.1	R	NA 5.2	UJ	NA 5.1	U	NA 5.1	U	NA 5.1	U	NA 5.5	U	NA 5.1	U
106-44-5 91-94-1	3&4-METHYLPHENOL 3.3-DICHLOROBENZIDINE	ug/l ug/l	5.1 5.1	R	5.1 5.1	R U	5.2 5.2	UJ	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U U*
101-55-3 106-47-8	4-BROMOPHENYL PHENYL ETHER 4-CHLOROANILINE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U U*
100-02-7	4-NITROPHENOL	ug/l	5.1	R	5.1	R	5.2	UJ	5.1	U	5.1	U	5.1	U	5.5	U	5.1	U
83-32-9 208-96-8	ACENAPHTHENE ACENAPHTHYLENE	ug/l ug/l	0.31	U	0.3	U	1 0.31	U	0.3	U	0.3	U	1 0.3	U	1.1 0.33	U	0.3	U
98-86-2 62-53-3	ACETOPHENONE ANILINE	ug/l ug/l	5.1 51	U	5.1 51	U	5.2 52	U	5.1 51	U	5.1 51	U	5.1 51	U	5.5 55	U	5.1 51	U
120-12-7	ANTHRACENE	ug/l	1	U	1	U	1	U	1	U	1	U	1	U	1.1	U	1	U
103-33-3 56-55-3	AZOBENZENE BENZO(A)ANTHRACENE	ug/l ug/l	5.1 0.31	U	5.1 0.3	U	5.2 0.31	U	5.1 0.3	U	5.1 0.3	U	5.1 0.3	U	5.5 0.33	U	5.1 0.3	U
50-32-8 205-99-2	BENZO(A)PYRENE BENZO(B)FLUORANTHENE	ug/l ug/l	0.2 0.31	U	0.2 0.3	U	0.21 0.31	U	0.2 0.3	U	0.2 0.3	U	0.2 0.3	U	0.22	U	0.2	U
191-24-2 207-08-9	BENZO(G,H,I)PERYLENE BENZO(K)FLUORANTHENE	ug/l	0.51 0.31	U	0.51 0.3	U	0.21 0.31	J	0.51 0.3	U	0.51 0.3	U	0.51 0.3	U	0.55 0.33	U	0.51 0.3	U
111-91-1	BIS(2-CHLOROETHOXY)METHANE	ug/l	5.1	U	5.1	U	5.2	U	5.1	U	5.1	U	5.1	U	5.5	U	5.1	U
111-44-4 108-60-1	BIS(2-CHLOROETHYL) ETHER BIS(2-CHLOROISOPROPYL)ETHER	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
117-81-7 85-68-7	BIS(2-ETHYLHEXYL) PHTHALATE BUTYLBENZYLPHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 0.34	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
218-01-9 53-70-3	CHRYSENE DIBENZ(A,H)ANTHRACENE	ug/l	1 0.51	U	1 0.51	U	1 0.42	U	1 0.51	U	1 0.51	U	1 0.51	U	1.1	U	1 0.51	U
132-64-9	DIBENZOFURAN	ug/l ug/l	5.1	U	5.1	U	5.2	U	5.1	U	5.1	U	5.1	U	5.5	U	5.1	U
84-66-2 131-11-3	DIETHYL PHTHALATE DIMETHYL PHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
84-74-2 117-84-0	DI-N-BUTYL PHTHALATE DI-N-OCTYL PHTHALATE	ug/l ug/l	5.1 5.1	U	5.1 5.1	U	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U	5.5 5.5	U	5.1 5.1	U
206-44-0	FLUORANTHENE	ug/l	0.37	J	0.25	J	1	U	1	U	1	U	1	U	1.1	U	1	U
86-73-7 118-74-1	FLUORENE HEXACHLOROBENZENE	ug/l ug/l	1	U	1	U	1	U	1	U	1	U	1	U	1.1 1.1	U	1	U
87-68-3 67-72-1	HEXACHLOROBUTADIENE HEXACHLOROETHANE	ug/l ug/l	0.61 5.1	U	0.61 5.1	U	0.62 5.2	U	0.61 5.1	U	0.61 5.1	U	0.61 5.1	U	0.66 5.5	U	0.61 5.1	U
193-39-5 78-59-1	INDENO(1,2,3-CD)PYRENE ISOPHORONE	ug/l ug/l	0.51 5.1	U	0.51 5.1	U	0.38 5.2	J	0.51 5.1	U	0.51 5.1	U	0.51 5.1	U	0.55 5.5	U	0.51 5.1	U
91-20-3	NAPHTHALENE	ug/l	1	U	1	U	1	U	1	U	1	U	1	U	1.1	U	1	U
98-95-3 106-44-5	P-CRESOL	ug/l ug/l	5.1 NA	U	5.1 NA	U	5.2 NA	U	5.1 NA	U	5.1 NA	U	5.1 NA	U	5.5 NA	U	5.1 NA	U
87-86-5 85-01-8	PENTACHLOROPHENOL PHENANTHRENE	ug/l ug/l	5.1 0.2	R	5.1 0.2	R U	5.2 0.21	UJ	5.1 0.2	U	5.1 0.2	U	5.1 0.2	U	5.5 0.22	U	5.1 0.2	U
108-95-2 129-00-0	PHENOL PYRENE	ug/l	5.1 0.28	R	5.1 5.1	R	5.2 5.2	UJ	5.1 5.1	UJ	5.1 5.1	UJ	5.1 5.1	UJ	5.5 5.5	UJ	5.1 5.1	UJ
35822-46-9	1,2,3,4,6,7,8-HpCDD	pg/l	NA	J	NA NA		NA		NA	U	NA	Ü	NA NA	Ü	NA	U	NA	Ĕ
67562-39-4 55673-89-7	1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
39227-28-6	1,2,3,4,7,8-HxCDD 1,2,3,4,7,8-HxCDD	pg/l	NA		NA		NA		NA		NA		NA		NA		NA	\perp
70648-26-9 57653-85-7	1,2,3,6,7,8-HxCDD	pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
57117-44-9 19408-74-3	1,2,3,6,7,8-HxCDF 1,2,3,7,8,9-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
72918-21-9 40321-76-4	1,2,3,7,8,9-HxCDF 1,2,3,7,8-PeCDD	pg/l pg/l	NA NA		NA NA	+	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	뻳
57117-41-6 60851-34-5	1,2,3,7,8-PeCDF 2,3,4,6,7,8-HxCDF	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	世
57117-31-4 1746-01-6	2,3,4,7,8-PeCDF 2,3,7,8-TCDD	pg/l pg/l	NA NA	L	NA NA	Ł	NA NA	Ł	NA NA	Ħ	NA NA		NA NA	Ħ	NA NA	F	NA NA	$oldsymbol{oldsymbol{eta}}$
51207-31-9 37871-00-4	2,3,7,8-TCDF HPCDD (TOTAL)	pg/l pg/l	NA NA	F	NA NA	Ł	NA NA	E	NA NA	Ħ	NA NA		NA NA	Ħ	NA NA		NA NA	$oldsymbol{\pm}$
38998-75-3 34465-46-8	HPCDF (TOTAL) HXCDD (TOTAL)	pg/l pg/l	NA NA	F	NA NA	╁	NA NA	Ŀ	NA NA	Ħ	NA NA		NA NA	Ħ	NA NA		NA NA	且
55684-94-1 3268-87-9	HXCDF (TOTAL) OCDD	pg/l pg/l	NA NA		NA NA	F	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	Ħ
39001-02-0 36088-22-9	OCDF PECDD (TOTAL)	pg/l pg/l	NA NA	F	NA NA	+	NA NA		NA NA	H	NA NA		NA NA	H	NA NA		NA NA	Ħ
30402-15-4 41903-57-5	PECDF (TOTAL) TCDD (TOTAL)	pg/l pg/l	NA NA	F	NA NA	+	NA NA		NA NA	H	NA NA		NA NA	H	NA NA		NA NA	Ħ
55722-27-5	TCDF (TOTAL)	pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
91-57-6 83-32-9	2-METHYLNAPHTHALENE ACENAPHTHENE	ug/l ug/l	1	U	1	U	1	U	NA NA	F	NA NA		NA NA	F	NA NA		NA NA	Ħ
208-96-8 120-12-7	ACENAPHTHYLENE ANTHRACENE	ug/l ug/l	0.3	U	0.31	U	0.31	U	NA NA		NA NA		NA NA		NA NA		NA NA	Ħ
56-55-3 50-32-8	BENZO(A)ANTHRACENE BENZO(A)PYRENE	ug/l ug/l	0.3 0.2	U	0.31 0.21	U	0.31 0.21	U	NA NA		NA NA		NA NA		NA NA		NA NA	丰
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3	U	0.31	U	0.31	U	NA		NA		NA		NA		NA	뻳
191-24-2 207-08-9	BENZO(G,H,I)PERYLENE BENZO(K)FLUORANTHENE	ug/l	0.4	U	0.31	U	0.42 0.31	U	NA NA	<u> </u>	NA NA		NA NA		NA NA		NA NA	뻳
EPH1122 EPH1122	C11-C22 AROMATICS, ADJUSTED C11-C22 AROMATICS, ADJUSTED	ug/l ug/l	290	J	460	J	100	U	100 NA	U	100 NA	U	100 NA	U	100 NA	U	NA	世
EPH1122 EPH1936	C11-C22 AROMATICS, UNADJUSTED C19-C36 ALIPHATICS	ug/l ug/l	290 100	J	470 130	J V	100 100	U	100 100	U	100 100	U	100 100	U	100 100	U	100 100	U
EPH1936 VPH912	C19-C36 ALIPHATICS (EPH-LL) C9-C18 ALIPHATICS	ug/l ug/l	100	U	100	U	100	U	NA 100	U	NA 100	U	NA 100	U	NA 100	U	100	U*
VPH912 218-01-9	C9-C18 ALIPHATICS (EPH-LL) CHRYSENE	ug/l ug/l	1	U	1	U	1	U	NA NA	É	NA NA		NA NA	É	NA NA	É	NA NA	Ħ
53-70-3 206-44-0	DIBENZ(A,H)ANTHRACENE FLUORANTHENE	ug/l ug/l	0.4 3.4	U		U	0.42	U	NA NA	H	NA NA		NA NA	H	NA NA	F	NA NA	丰
86-73-7 193-39-5	FLUORENE INDENO(1,2,3-CD)PYRENE	ug/l	1 0.4	U	1	U	1 0.42	U	NA NA NA	Ħ	NA NA NA		NA NA NA	Ħ	NA NA NA	F	NA NA NA	Ħ
91-20-3	NAPHTHALENE	ug/l	1	U	1	U	1	U	NA		NA		NA		NA		NA	$\downarrow \downarrow \downarrow$
85-01-8 129-00-0	PHENANTHRENE PYRENE	ug/l ug/l	0.62 1	J	0.37 1	J	0.21 1	U	NA NA		NA NA		NA NA		NA NA		NA NA	\pm
	TOTAL EPH	ug/l	290	J	600	J	100	U	100	U	100	כ	100	U	100	С	100	U

7440-36-0 7440-36-0 7440-38-2 7440-38-2 7440-39-3		epth:	6/25/2007 3 - 13 feet		6/25/2007		7/23/2007		12/11/2007		12/11/2007							R03-X
7440-36-0 7440-36-0 7440-38-2 7440-38-2	Analyte				3 - 13 feet		3 - 13 feet		4 - 14 feet		4.5 - 12.5 fee	t	12/11/2007 5 - 15 feet		12/11/2007 5 - 15 feet		12/11/200 3 - 13 feet	
7440-36-0 7440-38-2 7440-38-2		Units	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q
7440-38-2	ANTIMONY ANTIMONY (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA NA		NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-39-3	ARSENIC ARSENIC (DISSOLVED)	ug/l ug/l	NA 1.7	J	NA 1.7	J	NA NA		NA 1	U	NA 1	U	NA 10	V	NA 10	V	NA 1	U
7440-39-3	BARIUM (DISSOLVED)	ug/l ug/l	NA 27	V	NA 26	V	NA NA		NA 38	V	NA 11	V	NA 45	V	NA 44	V	NA 25	V
7440-41-7 7440-41-7	BERYLLIUM (DISSOLVED)	ug/l ug/l	NA 2	U	NA 1	U	NA NA		NA 2	U	NA 2	U	NA 2	U	NA 2	U	NA 2	U
7440-43-9 7440-43-9	CADMIUM (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA NA		NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7440-47-3 7440-47-3	CHROMIUM CHROMIUM (DISSOLVED)	ug/l ug/l	NA 9.8	V	NA 11	V	NA NA		NA 22	V	NA 15	V	NA 10	V	NA 9.1	V	NA 25	V
7440-50-8 7440-50-8	COPPER (DISSOLVED)	ug/l ug/l	NA NA		NA NA	Ė	NA NA		NA NA	Ť	NA NA	Ė	NA NA		NA NA		NA NA	İ
7439-92-1 7439-92-1	LEAD (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA NA		NA 1	U	NA 1	U	NA 1	U	NA 1	U	NA 1	U
7439-97-6 7439-97-6	MERCURY MERCURY (DISSOLVED)	ug/l ug/l	NA 0.2	U	NA 0.2	U	NA NA		NA 0.71	V	NA 0.2	U	NA 0.2	U	NA 0.2	U	NA 0.2	U
7440-02-0 7440-02-0	NICKEL (DISSOLVED)	ug/l ug/l	NA 3.6	V	NA 3.2	V	NA NA		NA 3.3	V	NA 2	U	NA 2	U	NA 2	U	NA 0.96	J
7782-49-2	SELENIUM	ug/l	NA		NA		NA		NA		NA		NA	Ŭ	NA		NA	
7782-49-2 7440-22-4	SELENIUM (DISSOLVED) SILVER	ug/l ug/l	2 NA	U	2 NA	U	NA NA		1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U
7440-22-4 7440-28-0	SILVER (DISSOLVED) THALLIUM	ug/l ug/l	1 NA	U	1 NA	U	NA NA		1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U
7440-28-0 7440-62-2	THALLIUM (DISSOLVED) VANADIUM	ug/l ug/l	1 NA	U	1 NA	U	NA NA		1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U
7440-62-2 7440-66-6	VANADIUM (DISSOLVED) ZINC	ug/l ug/l	0.91 NA	J	1 NA	J	NA NA		2 NA	U	2 NA	U	2 NA	U	2 NA	U	2 NA	U
7440-66-6	ZINC (DISSOLVED)	ug/l	9.4	٧	7.7	V	NA NA		2.4	J	1.9	J	1.5	J	1.2	J	20	V
630-20-6 71-55-6	1,1,1,2-TETRACHLOROETHANE 1,1,1-TRICHLOROETHANE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA	\Box	NA NA	Ш	NA NA	世
79-34-5 79-00-5	1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE	ug/l ug/l	2.5 5	U	2.5 5	U	NA NA		NA NA	Ы	NA NA		NA NA	oxed	NA NA	Ы	NA NA	± 1
75-34-3 75-35-4	1,1-DICHLOROETHANE 1,1-DICHLOROETHENE	ug/l ug/l	<u>5</u> 5	U	<u>5</u> 5	U	NA NA		NA NA	Н	NA NA	Ē	NA NA	ĿŢ	NA NA	H	NA NA	刊
563-58-6 87-61-6	1,1-DICHLOROPROPENE 1,2,3-TRICHLOROBENZENE	ug/l ug/l	<u>5</u> 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA	H	NA NA	H	NA NA	丑
96-18-4 120-82-1	1,2,3-TRICHLOROPROPANE 1,2,4-TRICHLOROBENZENE	ug/l ug/l	5 5	Ü	5 5	U	NA NA		NA NA		NA NA		NA NA	H	NA NA	H	NA NA	押
95-63-6 96-12-8	1,2,4-TRIMETHYLBENZENE 1,2-DIBROMO-3-CHLOROPROPANE	ug/l ug/l	5 25	Ü	5 25	U	NA NA		NA NA		NA NA		NA NA	H	NA NA		NA NA	丰
95-50-1 107-06-2	1,2-DICHLOROBENZENE 1,2-DICHLOROETHANE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\equiv
78-87-5 108678	1,2-DICHLOROPETHANE 1,2-DICHLOROPROPANE 1,3,5-TRIMETHYLBENZENE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA NA		NA NA		NA NA		NA NA	丰
541-73-1	1,3-DICHLOROBENZENE	ug/l	5	Ü	5	U	NA		NA		NA		NA		NA		NA	世
142-28-9 106-46-7	1,3-DICHLOROPROPANE 1,4-DICHLOROBENZENE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
123-91-1 594-20-7	1,4-DIOXANE 2,2-DICHLOROPROPANE	ug/l ug/l	250 5	U UJ	250 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
78-93-3 95-49-8	2-BUTANONE 2-CHLOROTOLUENE	ug/l ug/l	50 5	U	50 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
591-78-6 106-43-4	2-HEXANONE 4-CHLOROTOLUENE	ug/l ug/l	50 5	U	50 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
99-87-6 108-10-1	4-ISOPROPYLTOLUENE 4-METHYL-2-PENTANONE	ug/l ug/l	5 50	U	5 50	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	+
67-64-1 71-43-2	ACETONE BENZENE	ug/l ug/l	250 5	U	250 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	+
108-86-1 75-27-4	BROMOBENZENE BROMODICHLOROMETHANE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
75-25-2 74-83-9	BROMOFORM BROMOMETHANE	ug/l ug/l	5 10	U	5 10	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丰
75-15-0 56-23-5	CARBON DISULFIDE CARBON TETRACHLORIDE	ug/l ug/l	50 5	UJ U	50 5	UJ	NA NA		NA NA		NA NA		NA NA	H	NA NA		NA NA	丰
108-90-7 74-97-5	CHLOROBENZENE CHLOROBROMOMETHANE	ug/l ug/l	5	U	5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
75-00-3	CHLOROETHANE CHLOROFORM	ug/l ug/l	10 5	U	10	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
67-66-3 74-87-3	CHLOROMETHANE	ug/l	10	U	5 10	U	NA		NA		NA		NA		NA		NA	\pm
156-59-2 10061-01-5	CIS-1,2-DICHLOROETHENE CIS-1,3-DICHLOROPROPENE	ug/l ug/l	5 2.5	Ü	5 2.5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丗
124-48-1 74-95-3	DIBROMOCHLOROMETHANE DIBROMOMETHANE	ug/l ug/l	2.5 5	U	2.5 5	U	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
75-71-8 60-29-7	DICHLORODIFLUOROMETHANE DIETHYL ETHER	ug/l ug/l	5 50	U	5 50	U	NA NA		NA NA		NA NA		NA NA	\Box	NA NA	Ш	NA NA	世
108-20-3 100-41-4	DIISOPROPYL ETHER (DIPE) ETHYLBENZENE	ug/l ug/l	NA 5	U	NA 5	U	NA NA		NA NA	Ы	NA NA		NA NA	Ш	NA NA	Ы	NA NA	\pm
106-93-4 637-92-3	ETHYLENE DIBROMIDE ETHYLENE TERT-BUTYL ETHER	ug/l ug/l	5 25	U	5 25	U	NA NA		NA NA	Ы	NA NA	E	NA NA	oxdot	NA NA	H	NA NA	丑
87-68-3	HEXACHLOROBUTADIENE ISOPROPYL ETHER	ug/l ug/l	5 50	U	5 50	U	NA NA		NA NA	Ы	NA NA	Ē	NA NA	oxdot	NA NA	H	NA NA	丑
98-82-8	ISOPROPYLBENZENE M,P-XYLENES	ug/l ug/l	5 10	U	5 10	U	NA NA		NA NA	H	NA NA		NA NA	日	NA NA	H	NA NA	且
75-09-2 1634-04-4	METHYLENE CHLORIDE METHYL-T-BUTYL ETHER (MTBE)	ug/l ug/l	10 5	U	10	U	NA NA		NA NA	П	NA NA		NA NA	H	NA NA	П	NA NA	干
91-20-3 104-51-8	NAPHTHALENE N-BUTYLBENZENE	ug/l ug/l	25 5	U	25 5	U	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA	H	NA NA	干
103-65-1 95-47-6	N-PROPYLBENZENE O-XYLENE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA	Н	NA NA		NA NA	H	NA NA	H	NA NA	\blacksquare
135-98-8 100-42-5	SEC-BUTYLBENZENE STYRENE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA	H	NA NA		NA NA	Ħ	NA NA	Ħ	NA NA	#
994-05-8	TERT-AMYL METHYL ETHER (TAME) TERT-BUTYLBENZENE	ug/l	25	U	25	U	NA NA NA		NA NA NA	H	NA NA NA		NA NA NA	H	NA NA NA		NA NA	켇
98-06-6 127-18-4	TETRACHLOROETHYLENE	ug/l ug/l	5 5	Ū	5 5	U	NA		NA	H	NA		NA	Ħ	NA	H	NA	켇
109-99-9 108-88-3	TETRAHYDROFURAN TOLUENE	ug/l ug/l	50 5	U	50 5	U	NA NA		NA NA	H	NA NA		NA NA	Ħ	NA NA	H	NA NA	世
156-60-5 10061-02-6	TRANS-1,2-DICHLOROETHENE TRANS-1,3-DICHLOROPROPENE	ug/l ug/l	5 2.5	U	5 2.5	U	NA NA		NA NA	Ы	NA NA		NA NA	H	NA NA	H	NA NA	렏
79-01-6 75-69-4	TRICHLOROETHYLENE (TCE) TRICHLOROFLUOROMETHANE	ug/l ug/l	5 5	U	5 5	U	NA NA		NA NA		NA NA		NA NA		NA NA	Ы	NA NA	世
75-01-4	VINYL CHLORIDE	ug/l	5	U	5	U	NA		NA		NA		NA		NA		NA	+
71-43-2 VPH58	BENZENE C5-C8 ALIPHATICS, ADJUSTED	ug/l ug/l	5 100	U	5 100	U	NA NA		NA NA	Ы	NA NA	E	NA NA		NA NA	Ы	NA NA	$\pm \theta$
VPH58 VPH910	C5-C8 ALIPHATICS, UNADJUSTED C9-C10 AROMATICS	ug/l ug/l	100 100	U	100 100	U	NA NA		NA NA	Ħ	NA NA	F	NA NA	H	NA NA	H	NA NA	$\pm \overline{1}$
VPH912 VPH912	C9-C12 ALIPHATICS, ADJUSTED C9-C12 ALIPHATICS, UNADJUSTED	ug/l ug/l	100 100	U	100 100	U	NA NA		NA NA	H	NA NA	H	NA NA	H	NA NA	H	NA NA	$oldsymbol{+} \Box$
100-41-4	ETHYLBENZENE M.P-XYLENES	ug/l ug/l	5 10	U	5 10	U	NA NA		NA NA	H	NA NA		NA NA	H	NA NA	H	NA NA	丰
1634-04-4 91-20-3	METHYL-T-BUTYL ETHER (MTBE) NAPHTHALENE	ug/l ug/l	5 10	U	5 10	U	NA NA		NA NA	H	NA NA		NA NA	H	NA NA	H	NA NA	丰
95-47-6	O-XYLENE	ug/l ug/l	5	U	5 5	U	NA NA		NA NA		NA NA		NA NA	Ħ	NA NA	H	NA NA	丰
108-88-3	TOLUENE TOTAL VPH	ug/l ug/l	5 100	U	5 100	U	NA NA		NA NA	Н	NA NA	1	NA NA	+-+	NA NA	++	NA NA	+

Notes:
ug/L - micrograms per liter
V - Valid result, no qualification needed
U - not detected, laboratory reporting limit listed
J - concentration (or reporting limit) is estimated
DDA - Demolition Debris Area
NA - Not analyzed

	Sta Field Samp	ation:	DD-MW-201	•	DD-MW-201)6-D	DD-MW-002 DD-MW-002-R02	2-X	DD-MW-203 DD-MW-203-R06	-X	DD-MW-204 DD-MW-204-R0	5-X	DD-MW-205 DD-MW-205-R0	5-X	DD-MW-206		DD-MW-207 DD-MW-207-R0	6-X	DD-MW-208	-
	Sample		5/19/2008 4 - 14 feet	00 X	5/19/2008 4 - 14 feet	,	5/19/2008 9.8 - 14.8 feet		5/19/2008 4.5 - 12.5 feet	Α	5/19/2008 7.5 - 13.5 feet		5/19/2008 5 - 11 feet	<i>0</i>	5/19/2008 5 - 11 feet	,0 X	5/19/2008 5 - 15 feet	0 /	5/19/2008 3 - 13 feet	
CAS Number		Units	Result	Q	Result	Q		Q		Q	Result	Q		Q	Result	Q		Q	Result	Q
120-82-1 95-50-1	1,2,4-TRICHLOROBENZENE 1,2-DICHLOROBENZENE	ug/l ug/l	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U	5.1	U	5.1 5.1	U		U	5.1 5.1	U	5.2	U	5.1 5.1	U
541-73-1 106-46-7	1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE	ug/l ug/l	5.2 5.2	U	5.1 5.1	U		U	5.1	U	5.1 5.1	U	***	U	5.1 5.1	U	5.2	U	5.1 5.1	U
95-95-4 88-06-2	2,4,5-TRICHLOROPHENOL 2,4,6-TRICHLOROPHENOL	ug/l ug/l	5.2 5.2	UJ	5.1 5.1	UJ		UJ		U UJ	5.1 5.1	UJ		UJ		UJ		U	5.1 5.1	U
120-83-2 105-67-9	2,4-DICHLOROPHENOL 2,4-DIMETHYLPHENOL	ug/l ug/l	5.2 5.2	UJ	5.1 5.1	UJ		UJ		UJ	5.1 5.1	UJ		UJ		UJ		U	5.1 5.1	U
51-28-5 121-14-2	2,4-DINITROPHENOL 2.4-DINITROTOLUENE	ug/l ug/l	5.2 5.2	UJ	5.1 5.1	UJ		UJ	5.1	UJ	0.84 5.1	J	5.1 5.1	UJ		UJ	5.2	U	5.1 5.1	U
606-20-2	2,6-DINITROTOLUENE	ug/l	5.2	U	5.1	U	5.1	U	5.1	U	5.1	U	5.1	U	5.1	U	5.2	U	5.1	U
91-58-7 95-57-8	2-CHLORONAPHTHALENE 2-CHLOROPHENOL	ug/l ug/l	5.2 5.2	U	5.1 5.1	U		UJ	5.1	UJ	5.1 5.1	U UJ	5.1	UJ		UJ	5.2	U	5.1 5.1	U
91-57-6 95-48-2	2-METHYLNAPHTHALENE 2-METHYLPHENOL	ug/l ug/l	1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U	1 NA	U
95-48-7 88-75-5	2-METHYLPHENOL (O-CRESOL) 2-NITROPHENOL	ug/l ug/l	5.2 5.2	UJ	5.1 5.1	IJ	5.1 5.1	UJ		UJ	5.1 5.1	UJ	_	UJ	5.1 5.1	UJ		U	5.1 5.1	U
106-44-5 91-94-1	3&4-METHYLPHENOL 3,3-DICHLOROBENZIDINE	ug/l ug/l	NA 5.2	U	NA 5.1	U	NA 5.1	U	NA	U	NA 5.1	U	NA 5.1	U	NA 5.1	U	NA	U	NA 5.1	U
101-55-3	4-BROMOPHENYL PHENYL ETHER 4-CHLOROANILINE	ug/l	5.2	Ü	5.1	U	5.1	U	5.1	U	5.1 5.1	U	5.1	U	5.1	U	5.2	U	5.1	U
106-47-8 100-02-7	4-NITROPHENOL	ug/l ug/l	5.2 5.2	U	5.1 5.1	U		UJ	5.1	UJ	5.1	UJ		UJ	5.1 5.1	UJ	5.2	UJ	5.1 5.1	UJ
83-32-9 208-96-8	ACENAPHTHENE ACENAPHTHYLENE	ug/l ug/l	0.31	U	0.31	U	0.3	U		U	0.3	U	0.3	U	0.3	U		U	0.3	U
98-86-2 62-53-3	ACETOPHENONE ANILINE	ug/l ug/l	5.2 52	U	5.1 51	U	5.1 51	U		U UJ	5.1 51	U	5.1 51	U	5.1 51	U		U	5.1 51	U
120-12-7 103-33-3	ANTHRACENE AZOBENZENE	ug/l ug/l	1 5.2	U	1 5.1	U	1 5.1	U	1	U	1 5.1	U	1	U	1 5.1	U	1	U	1 5.1	U
56-55-3 50-32-8	BENZO(A)ANTHRACENE BENZO(A)PYRENE	ug/l ug/l	0.31 0.21	U	0.31 0.2	U	0.3	U	0.3	U	0.3	U	_	U	0.3	U	0.31	U	0.3	U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.31	U	0.31	U	0.3	U	0.3	U	0.3	U	0.3	U	0.3	U	0.31	U	0.3	U
191-24-2 207-08-9	BENZO(G,H,I)PERYLENE BENZO(K)FLUORANTHENE	ug/l ug/l	0.52 0.31	U	0.51 0.31	U	0.51 0.3	U	0.3	U	0.51 0.3	U	0.3	U	0.51 0.3	U	0.31	U	0.51 0.3	U
111-91-1 111-44-4	BIS(2-CHLOROETHOXY)METHANE BIS(2-CHLOROETHYL) ETHER	ug/l ug/l	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U		U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U		U	5.1 5.1	U
108-60-1 117-81-7	BIS(2-CHLOROISOPROPYL)ETHER BIS(2-ETHYLHEXYL) PHTHALATE	ug/l ug/l	5.2 0.48	U	5.1 5.1	U	5.1	U	5.1	U	5.1 0.45	U	5.1	U	5.1 0.49	U	5.2	U	5.1 1.6	U
85-68-7 218-01-9	BUTYLBENZYLPHTHALATE CHRYSENE	ug/l	5.2	U	5.1	U	5.1	U	5.1	U	5.1	U	5.1	U	5.1	U	5.2	U	5.1	U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l ug/l	0.52	Ü	0.51	U	0.51	U	0.51	U	0.51	Ü	0.51	U	0.51	U	0.52	U	0.51	U
132-64-9 84-66-2	DIBENZOFURAN DIETHYL PHTHALATE	ug/l ug/l	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U		U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U		U	5.1 5.1	U
131-11-3 84-74-2	DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE	ug/l ug/l	5.2 5.2	U	5.1 5.1	U	5.1 5.1	U		U	5.1 5.1	U	5.1 5.1	U	5.1 5.1	U		U	5.1 1.5	U
117-84-0 206-44-0	DI-N-OCTYL PHTHALATE FLUORANTHENE	ug/l ug/l	5.2	U	5.1	U	5.1	U	5.1	U	5.1	U		U	5.1	U	5.2	U	5.1	U
86-73-7	FLUORENE	ug/l	1	Ū	1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U
118-74-1 87-68-3	HEXACHLOROBENZENE HEXACHLOROBUTADIENE	ug/l ug/l	1 0.62	U	1 0.61	U	1 0.61	U	0.61	U	1 0.61	U		U	0.61	U	0.62	U	0.61	U
67-72-1 193-39-5	HEXACHLOROETHANE INDENO(1,2,3-CD)PYRENE	ug/l ug/l	5.2 0.52	U	5.1 0.51	U	5.1 0.51	U		U	5.1 0.51	U		U	5.1 0.51	U	0.52	U	5.1 0.51	U
78-59-1 91-20-3	ISOPHORONE NAPHTHALENE	ug/l ug/l	5.2 1	U	5.1 1	\Box	5.1 1	U		U	5.1 1	U	_	U	5.1 1	U		U	5.1 1	U
98-95-3 106-44-5	NITROBENZENE P-CRESOL	ug/l ug/l	5.2 5.2	U	5.1 5.1	U		U UJ	5.1	U	5.1 5.1	U	_	U	5.1 5.1	U	5.2	U	5.1 5.1	U
87-86-5	PENTACHLOROPHENOL	ug/l	1	UJ	1	UJ	1	UJ	1	UJ	1	UJ	1	UJ	1	UJ	1	U	1	U
85-01-8 108-95-2	PHENANTHRENE PHENOL	ug/l ug/l	0.21 5.2	U	0.2 5.1	U		U	5.1	UJ	0.2 5.1	UJ	5.1	UJ	5.1	UJ	5.2	U	0.2 5.1	UJ
129-00-0	PYRENE	ug/l	5.2	U	5.1	U	5.1	U		U	5.1	U	<u> </u>	U	<u> </u>	U	Ų	U	5.1	U
35822-46-9 67562-39-4	1,2,3,4,6,7,8-HpCDD 1,2,3,4,6,7,8-HpCDF	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
55673-89-7 39227-28-6	1,2,3,4,7,8,9-HpCDF 1,2,3,4,7,8-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
70648-26-9 57653-85-7 57117-44-9	1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDD 1,2,3,6,7,8-HxCDF	pg/l pg/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	\pm
19408-74-3	1,2,3,7,8,9-HxCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA NA		NA NA		NA NA		NA NA	\pm
72918-21-9 40321-76-4	1,2,3,7,8,9-HxCDF 1,2,3,7,8-PeCDD	pg/l pg/l	NA NA NA		NA		NA		NA		NA		NA		NA		NA		NA	\pm
57117-41-6 60851-34-5	1,2,3,7,8-PeCDF 2,3,4,6,7,8-HxCDF	pg/l pg/l	NA		NA NA		NA NA		NA NA		NA NA NA		NA NA		NA NA NA		NA NA NA		NA NA NA	\pm
57117-31-4 1746-01-6 51207-31-9	2,3,4,7,8-PeCDF 2,3,7,8-TCDD	pg/l pg/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	\pm
37871-00-4	2,3,7,8-TCDF HPCDD (TOTAL)	pg/l pg/l	NA NA NA		NA		NA		NA		NA		NA		NA		NA		NA	\pm
38998-75-3 34465-46-8	HYCDD (TOTAL)	pg/l pg/l	NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
55684-94-1 3268-87-9	HXCDF (TOTAL) OCDD	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
39001-02-0 36088-22-9	PECDD (TOTAL)	pg/l pg/l	NA NA		NA NA		NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA	\pm
30402-15-4 41903-57-5	PECDF (TOTAL) TCDD (TOTAL)	pg/l pg/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
55722-27-5 91-57-6	TCDF (TOTAL) 2-METHYLNAPHTHALENE	pg/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	+
83-32-9 208-96-8	ACENAPHTHENE ACENAPHTHYLENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	Ħ
120-12-7	ANTHRACENE BENZO(A)ANTHRACENE	ug/l ug/l	NA NA		NA NA		NA NA NA		NA NA NA		NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA	Ħ
56-55-3 50-32-8 205-99-2	BENZO(A)ANTHRACENE BENZO(A)PYRENE BENZO(B)FLUORANTHENE	ug/l ug/l	NA NA NA	+	NA NA NA		NA NA NA		NA NA NA	#	NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA	\mp
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	NA NA NA		NA		NA	Ħ	NA	1	NA NA NA		NA NA NA		NA NA NA		NA NA NA	H	NA NA NA	\pm
207-08-9 EPH1122	BENZO(K)FLUORANTHENE C11-C22 AROMATICS, ADJUSTED	ug/l ug/l	NA		NA NA	ļ.,	NA NA		NA NA		NA		NA		NA		NA	,,	100	U
EPH1122 EPH1122 EPH1936	C11-C22 AROMATICS, ADJUSTED C11-C22 AROMATICS, UNADJUSTED		100 100	U	100 100	U	100 100	U	100	U	100 100	U		U	100 100	U	100	U	100	U
EPH1936	C19-C36 ALIPHATICS C19-C36 ALIPHATICS (EPH-LL)	ug/l ug/l	NA 100	U	NA 100	U	NA 100	U		U	NA 100	U		U	NA 100	U		U	100	
VPH912 VPH912	C9-C18 ALIPHATICS C9-C18 ALIPHATICS (EPH-LL)	ug/l ug/l	NA 100	U	NA 100	U	NA 100	U		U	NA 100	U		U		U		U	100	U
218-01-9 53-70-3	CHRYSENE DIBENZ(A,H)ANTHRACENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	#	NA NA		NA NA		NA NA		NA NA	H	NA NA	\ddagger
206-44-0 86-73-7	FLUORANTHENE FLUORENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA	\pm
193-39-5 91-20-3	INDENO(1,2,3-CD)PYRENE NAPHTHALENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA	Ы	NA NA	峀
85-01-8 129-00-0	PHENANTHRENE PYRENE	ug/l ug/l	NA NA	ŧ.	NA NA		NA NA	Ы	NA NA	1	NA NA		NA NA		NA NA	L	NA NA		NA NA	Ħ
	TOTAL EPH	ug/l	100	U	100	U	100	U	100	U	100	U	100	U	100	U	100	U	100	U

	Sta Field Samp	ation:	DD-MW-201-R0		DD-MW-201 DD-MW-201-R0	6-D	DD-MW-002 DD-MW-002-R02	2-X	DD-MW-203 DD-MW-203-R06	S-X	DD-MW-204 DD-MW-204-R05	5-X	DD-MW-205 DD-MW-205-R05	5-X	DD-MW-206-R0		DD-MW-207	6-X	DD-MW-20 DD-MW-208-R	_
		epth:	5/19/2008 4 - 14 feet		5/19/2008 4 - 14 feet		5/19/2008 9.8 - 14.8 feet		5/19/2008 4.5 - 12.5 feet		5/19/2008 7.5 - 13.5 feet		5/19/2008 5 - 11 feet		5/19/2008 5 - 11 feet		5/19/2008 5 - 15 feet		5/19/2008 3 - 13 feet	i
7440-36-0	Analyte ANTIMONY	Units ug/l	Result NA	Q	Result NA	Q	Result NA	Q	Result NA	Q	Result NA	Q	Result	Q	Result NA	Q	Result NA	Q	Result	Q
7440-36-0 7440-38-2	ANTIMONY (DISSOLVED) ARSENIC	ug/l ug/l	1 NA	U	1 NA	U	NA	U	1 NA	UJ	1 NA	U	NA	U	1 NA	U	1 NA	U	1 NA	U
7440-38-2 7440-39-3	ARSENIC (DISSOLVED) BARIUM	ug/l ug/l	1 NA	U	1 NA	U	NA	U	1 NA	U	2 NA	U	NA	U	1 NA	U	3.9 NA	V	1 NA	U
7440-39-3 7440-41-7	BARIUM (DISSOLVED) BERYLLIUM	ug/l ug/l	46 NA	V	45 NA	V	NA	V	11 NA	V	20 NA	V	NA	V	NA	V	31 NA	V	31 NA	V
7440-41-7 7440-43-9 7440-43-9	BERYLLIUM (DISSOLVED) CADMIUM CADMIUM (DISSOLVED)	ug/l ug/l ug/l	1 NA 1	U	1 NA 1	U	NA	U	1 NA 1	U	1 NA 1	U	NA NA	U	NA	U	1 NA 1	U	1 NA 1	U
7440-47-3 7440-47-3	CHROMIUM CHROMIUM (DISSOLVED)	ug/l ug/l	NA 2.1	J	NA 3.5	J	NA	V	NA 0.63	J	NA 1	IJ	NA		NA 0.51	J.	NA 1.3	٧	NA 0.36	
7440-50-8 7440-50-8	COPPER (DISSOLVED)	ug/l ug/l	NA NA		NA NA	J	NA NA	Ť	NA NA	Ů	NA NA	_	NA NA		NA NA	Ů	NA NA	_	NA NA	Ť
7439-92-1 7439-92-1	LEAD LEAD (DISSOLVED)	ug/l ug/l	NA 1	U	NA 1	U	NA	U	NA 1	U	NA 1	U	NA	U	NA	U	NA 1	U	NA 0.46	J
7439-97-6 7439-97-6	MERCURY MERCURY (DISSOLVED)	ug/l ug/l	NA 0.2	U	NA 0.2	U		U	NA 0.2	U	NA 0.2	U	·	U	NA 0.2	U	NA 0.2	U	NA 0.2	U
7440-02-0 7440-02-0	NICKEL NICKEL (DISSOLVED)	ug/l ug/l	NA 1.3	V	NA 1.2	V		V	NA 1.2	V	NA 1.2	V		J	NA 0.69	J	NA 0.43	J	NA 2	V
7782-49-2 7782-49-2	SELENIUM SELENIUM (DISSOLVED) SILVER	ug/l	NA 6.8	V	NA 7.8	V		U		U	NA 2	U	NA 1 NA	U	NA 1 NA	U	NA 1 NA	U	NA 1	U
7440-22-4 7440-22-4 7440-28-0	SILVER SILVER (DISSOLVED) THALLIUM	ug/l ug/l ug/l	NA 1 NA	U	NA 1 NA	U	NA 1 NA	U	NA 1 NA	U	NA 1 NA	U		U		U	1 NA	U	NA 1 NA	U
7440-28-0 7440-28-0 7440-62-2	THALLIUM (DISSOLVED) VANADIUM	ug/I ug/I	1 NA	U	1 NA	U		U	1 NA	U	1 NA	U		U		U		U	1 NA	U
7440-62-2 7440-66-6	VANADIUM (DISSOLVED) ZINC	ug/l ug/l	1 NA	U	0.18 NA	J		U	0.25 NA	J	1 NA	U		J	1 NA	U	1 NA	U	0.52 NA	J
7440-66-6	ZINC (DISSOLVED)	ug/l	2.7	V	2.3	J	1.2	J	5.3	V	5	U	1.6	J	1.9	J	1.9	J	110	V
630-20-6 71-55-6 79-34-5	1,1,1,2-TETRACHLOROETHANE 1,1,1-TRICHLOROETHANE	ug/l ug/l	NA NA	H	NA NA NA		NA NA NA	1	NA NA NA	4	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA	丗
79-34-5 79-00-5 75-34-3	1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE 1,1-DICHLOROETHANE	ug/l ug/l ug/l	NA NA NA	Ħ	NA NA NA		NA NA NA	1	NA NA NA	4	NA NA NA		NA NA NA		NA NA NA	Ļ	NA NA NA		NA NA NA	텎
75-34-3 75-35-4 563-58-6	1,1-DICHLOROETHANE 1,1-DICHLOROETHENE 1,1-DICHLOROPROPENE	ug/I ug/I	NA NA NA	Ħ	NA NA NA		NA NA	1	NA NA NA	4	NA NA		NA NA		NA NA NA	F	NA NA NA		NA NA NA	丰
87-61-6 96-18-4	1,2,3-TRICHLOROBENZENE 1,2,3-TRICHLOROPROPANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
120-82-1 95-63-6	1,2,4-TRICHLOROBENZENE 1,2,4-TRIMETHYLBENZENE	ug/l ug/l	NA NA	Е	NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA	E	NA NA		NA NA	田
96-12-8 95-50-1	1,2-DIBROMO-3-CHLOROPROPANE 1,2-DICHLOROBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
107-06-2 78-87-5	1,2-DICHLOROETHANE 1,2-DICHLOROPROPANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
108678 541-73-1 142-28-9	1,3,5-TRIMETHYLBENZENE 1,3-DICHLOROBENZENE 1,3-DICHLOROPROPANE	ug/l ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	〓
106-46-7 123-91-1	1,4-DICHLOROBENZENE 1,4-DIOXANE	ug/I ug/I	NA NA 3.1	UJ	NA NA 3.1	UJ	NA	II.I	NA NA 3	IJJ	NA	UJ	NA	IJJ	NA	UJ	NA NA 3.1	IJJ	NA NA 3	UJ
594-20-7 78-93-3	2,2-DICHLOROPROPANE 2-BUTANONE	ug/l ug/l	NA NA	00	NA NA	00	NA NA	00	NA NA	00	NA NA	00	NA NA	00	NA NA	0.0	NA NA	00	NA NA	
95-49-8 591-78-6	2-CHLOROTOLUENE 2-HEXANONE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
106-43-4 99-87-6	4-CHLOROTOLUENE 4-ISOPROPYLTOLUENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
108-10-1 67-64-1	4-METHYL-2-PENTANONE ACETONE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
71-43-2 108-86-1	BENZENE BROMOBENZENE BROMODICHLOROMETHANE	ug/l ug/l	NA NA	L	NA NA		NA NA	=	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
75-27-4 75-25-2 74-83-9	BROMODICHLOROME I HANE BROMOFORM BROMOMETHANE	ug/l ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	\pm
75-15-0 56-23-5	CARBON DISULFIDE CARBON TETRACHLORIDE	ug/l ug/l	NA NA	H	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丰
108-90-7 74-97-5	CHLOROBENZENE CHLOROBROMOMETHANE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\blacksquare
75-00-3 67-66-3	CHLOROETHANE CHLOROFORM	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	
74-87-3 156-59-2	CHLOROMETHANE CIS-1,2-DICHLOROETHENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\pm
10061-01-5 124-48-1 74-95-3	CIS-1,3-DICHLOROPROPENE DIBROMOCHLOROMETHANE DIBROMOMETHANE	ug/l ug/l ug/l	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	\pm
74-95-3 75-71-8 60-29-7	DICHLORODIFLUOROMETHANE DIETHYL ETHER	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA NA		NA NA		NA NA		NA NA		NA NA	\mp
108-20-3 100-41-4	DIISOPROPYL ETHER (DIPE) ETHYLBENZENE	ug/l ug/l	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	\mp
106-93-4 637-92-3	ETHYLENE DIBROMIDE ETHYLENE TERT-BUTYL ETHER	ug/l ug/l	NA NA	H	NA NA		NA NA		NA NA		NA NA	E	NA NA	E	NA NA	F	NA NA		NA NA	H
87-68-3	HEXACHLOROBUTADIENE ISOPROPYL ETHER	ug/l ug/l	NA NA	Ы	NA NA		NA NA	\exists	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	目
98-82-8	ISOPROPYLBENZENE M,P-XYLENES	ug/l ug/l	NA NA	H	NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丗
75-09-2 1634-04-4	METHYLENE CHLORIDE METHYL-T-BUTYL ETHER (MTBE)	ug/l ug/l	NA NA NA	H	NA NA NA		NA NA NA	1	NA NA NA	\dashv	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	\mp
91-20-3 104-51-8 103-65-1	NAPHTHALENE N-BUTYLBENZENE N-PROPYLBENZENE	ug/l ug/l ug/l	NA NA NA	H	NA NA NA		NA NA NA	1	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	丰
95-47-6 135-98-8	O-XYLENE SEC-BUTYLBENZENE	ug/l ug/l	NA NA	H	NA NA	H	NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丰
100-42-5 994-05-8	STYRENE TERT-AMYL METHYL ETHER (TAME)	ug/l ug/l	NA NA	H	NA NA		NA NA]	NA NA		NA NA		NA NA		NA NA	L	NA NA		NA NA	丑
98-06-6 127-18-4	TERT-BUTYLBENZENE TETRACHLOROETHYLENE	ug/l ug/l	NA NA	F	NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丑
109-99-9 108-88-3	TETRAHYDROFURAN TOLUENE	ug/l ug/l	NA NA	H	NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA		NA NA	丗
156-60-5 10061-02-6 79-01-6	TRANS-1,2-DICHLOROETHENE TRANS-1,3-DICHLOROPROPENE TRICHLOROETHYLENE (TCE)	ug/l ug/l ug/l	NA NA NA	H	NA NA NA		NA NA NA	1	NA NA NA	4	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	井
79-01-6 75-69-4 75-01-4	TRICHLOROETHYLENE (TCE) TRICHLOROFLUOROMETHANE VINYL CHLORIDE	ug/l ug/l	NA NA NA	H	NA NA NA		NA NA NA	#	NA NA NA	4	NA NA NA		NA NA NA		NA NA NA		NA NA NA		NA NA NA	丰
71-43-2	BENZENE	ug/I	NA NA		NA NA		NA NA		NA NA	4	NA NA		NA NA		NA NA	F	NA NA		NA NA	肀
VPH58 VPH58	C5-C8 ALIPHATICS, ADJUSTED C5-C8 ALIPHATICS, UNADJUSTED	ug/l ug/l	NA NA		NA NA		NA NA]	NA NA		NA NA	E	NA NA	E	NA NA	E	NA NA		NA NA	
VPH910 VPH912	C9-C10 AROMATICS C9-C12 ALIPHATICS, ADJUSTED	ug/l ug/l	NA NA	Ħ	NA NA		NA NA	1	NA NA	1	NA NA		NA NA		NA NA		NA NA		NA NA	坦
VPH912 100-41-4	C9-C12 ALIPHATICS, UNADJUSTED ETHYLBENZENE	ug/l	NA NA	H	NA NA		NA NA	1	NA NA		NA NA		NA NA		NA NA	L	NA NA		NA NA	\pm
1634-04-4 91-20-3	M,P-XYLENES METHYL-T-BUTYL ETHER (MTBE) NAPHTHALENE	ug/l ug/l ug/l	NA NA NA	H	NA NA NA		NA NA NA	1	NA NA NA	4	NA NA NA		NA NA NA		NA NA NA	L	NA NA NA		NA NA NA	‡
91-20-3 95-47-6 108-88-3	O-XYLENE TOLUENE	ug/I ug/I	NA NA NA	H	NA NA NA		NA NA	7	NA NA NA	4	NA NA NA		NA NA		NA NA NA	F	NA NA NA		NA NA NA	丰
	TOTAL VPH	ug/l	NA NA		NA NA		NA NA		NA NA		NA		NA NA		NA NA		NA NA		NA NA	工

Notes:
ug/L - micrograms per liter
V - Valid result, no qualification needed
U - not detected, laboratory reporting limit listed
J - concentration (or reporting limit) is estimated
DDA - Demolition Debris Area
NA - Not analyzed

Table 6-1 **Groundwater Quality** Phase II CSA Addendum **Demolition Debris Area** Bird Machine Company - Walpole, MA

					Upgradier	nt A	Area of DDA						Inside DDA						Downg	gradie	ent of DDA			_
	Si	tation:		DD-M	IW-204	T	DD)-M\	W-205	T	DD-MW-001		D	D-M	1W-002	T			D	D-MV	V-201			
	Field Sam	ple ID:	DD-MW-204-R0	02-X	DD-MW-204-R05-2	Х	DD-MW-205-R03-	-X	DD-MW-205-R05-	Х	DD-MW-001-R02	2-X	DD-MW-002-R0)1-X	DD-MW-002-R02-2	Х	DD-MW-201-R04	4-X	DD-MW-201-R0	5-X	DD-MW-201-R06	S-X	DD-MW-201-R06	j-D
	Sample	Date:	6/5/2007		5/19/2008		6/6/2007		5/19/2008		6/5/2007		6/5/2007		5/19/2008		6/5/2007		12/11/2007		5/19/2008		5/19/2008	
		Depth:	7.5 - 13.5 fee	et	7.5 - 13.5 feet		5 - 11 feet		5 - 11 feet		11.6 - 16.6 fee	t	9.8 - 14.8 fee	et	9.8 - 14.8 feet		4 - 14 feet		4 - 14 feet	Ī	4 - 14 feet		4 - 14 feet	
CAS Number	Analyte	Únits	Result	Q	Result (Q	Result	Q	Result	Q	Result	Q	Result	Q	Result (Q	Result	Q	Result	Q	Result	Q	Result	Q
51-28-5	2,4-DINITROPHENOL	ug/l	5.1	UJ	0.84	J	5.1	UJ	5.1 l	IJ	5.1	UJ	5.1	UJ	5.1 L	JJ	5.3	UJ	5.1	UJ	5.2	UJ	5.1	UJ
56-55-3	BENZO(A)ANTHRACENE	ug/l		U		Ü	0.3	U		U	0.3	U	0.3	U		U	0.32	U	0.3	U		U		U
50-32-8	BENZO(A)PYRENE	ug/l		U		Ü		Ū		Ū	0.2	Ü	0.2	Ü		Ü	0.21	Ü	0.2	Ü		Ü	0.2	Ü
205-99-2	BENZO(B)FLUORANTHENE	ug/l		Ü	***	Ü		U		Ü	0.3	Ü	0.3	Ü		U	0.32	U	0.3	Ü	*	Ü	-	Ü
191-24-2	BENZO(G,H,I)PERYLENE	ug/l		U		Ü		U		Ü	0.51	U	0.51	Ü		U	0.53	U	0.51	Ü		U		U
207-08-9	BENZO(K)FLUORANTHENE	ug/l		U		Ü		U		Ü	0.3	U	0.3	U		U	0.32	U	0.3	Ü		Ü		Ü
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l		Ü		J		U		U	5.1	U	5.1	U		U	5.3	U	5.1	Ü		J	5.1	U
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.1	U		U		U		U	5.1	U	5.1	U		U	5.3	U	5.1	U		U	5.1	U
	CHRYSENE		1	U		U		U		U	J. I	U	1	U		U		U	1	U		U	1	U
218-01-9 53-70-3	DIBENZ(A.H)ANTHRACENE	ug/l ug/l	0.51	U		U		U		U	0.51	U	0.51	U		U	1.1 0.53	U	0.51	U		U	0.51	U
	, ,					_		_		_						_						-		
84-74-2	DI-N-BUTYL PHTHALATE	ug/l		U	5.1 l	_		U		U	5.1	U	5.1	U		U	5.3	U	5.1	U		U		U
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	_	U		U		U		U		U	5.1	U		U	5.3	U		U		U		U
206-44-0	FLUORANTHENE	ug/l		U		U		U		U		U	1	U		U	1.1	U		U		U		U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.51	U	0.51 l	_		U		U		U	0.51	U		U	0.53	U		U		U		U
85-01-8	PHENANTHRENE	ug/l	0.2	U	0.2 l	_		U		U	0.2	U	0.2	U		U	0.21	U		U		U	0.2	U
129-00-0	PYRENE	ug/l	5.1	U	5.1 l	U	5.1	U	5.1	U	5.1	U	5.1	U	5.1 l	U	5.3	J	5.1	U	5.2	U	5.1	U
56-55-3	BENZO(A)ANTHRACENE	ua/l	0.3	U	NA	7	0.3	U	NA	-	0.3	U	0.32	U	NA	-	0.32	U	NA	-	NA	_	NA	
50-32-8	BENZO(A)PYRENE	ug/I		U	NA NA	+		IJ	NA NA	-	0.2	U	0.32	IJ		\dashv	0.32	U	NA NA		NA NA		NA NA	-
205-99-2	BENZO(B)FLUORANTHENE	ug/I		U		_	*	U	NA NA	十		U	0.32	Ü		-	0.32	U	NA NA		NA NA		NA NA	-
191-24-2	BENZO(G.H.I)PERYLENE	ug/l		Ü	NA NA	T		IJ	NA NA	7	0.4	U	0.42	IJ		1	0.42	IJ	NA NA	t	NA NA		NA NA	-
207-08-9	BENZO(K)FLUORANTHENE	ua/l		Ŭ			• • •	Ü	NA		0.3	Ŭ	0.32	Ŭ			0.32	Ŭ	NA NA		NA		NA	
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	NA		100	U	NA		100	Ū	110	U	NA		110	U	100	U	NA		NA	
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	NA		100	U	NA		100	U	110	U	NA		110	כ	100	U	NA		NA	
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.4	U	NA		0.4	U	NA		0.4	U	0.42	U			0.42	J	NA		NA		NA	
206-44-0	FLUORANTHENE	ug/l		U				U	NA			U	1.1	U			1.1	U	NA		NA		NA	
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l		U	NA		***	U	NA		0.4	U	0.42	U			0.42	U	NA		NA		NA	
85-01-8	PHENANTHRENE	ug/l	0.2	U	NA	_	0.2	U	NA	_	0.2	U	0.21	U	NA	_	0.21	U	NA	_	NA	_	NA	_
7440-36-0	ANTIMONY (DISSOLVED)	ua/l	1	U	1 1	U	1	U	1	U	1	U	1	U	1 1	U	1	U	1	U	1	U	1	U
7440-36-0	ARSENIC (DISSOLVED)	ug/I		IJ		U		U		U	2	U	2	IJ		U	0.81	.]	1	IJ		U	1	U
7440-39-3	BARIUM (DISSOLVED)	ug/I		Ť	20	~	18	Ť	7.6	┪	47		32	T	19	1	44	Ť	38	- 	46	-	45	\dashv
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l		U		U		U		U		U	1	U		U	1	U		U		U		U
7440-43-9	CADMIUM (DISSOLVED)	ug/l		Ü	1 1	Ü	1	Ū	1	Ü	1	Ü	1	Ü	1 1	U	1	د (1	Ü	1	Ŭ		Ü
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	11		1 l	U	8.4		0.52	J	35		7.3		1.9		25		22		2.1	J	3.5	J
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA		NA		NA		NA		NA		NA		NA		NA		NA	
7439-92-1	LEAD (DISSOLVED)	ug/l		U		U		U		U	1	U	1	U		U	1	J	1	U		U	1	U
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U		U	0.2	U		U	0.2	U	0.2	U	0.2	U	0.2	כ	0.71	[V.E	U	0.2	U
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.1	1	1.2	_	1.4	ᆜ	****	J	7.2		2.5	L	1.2	_	1.3	<u> </u>	3.3		1.3		1.2	
7782-49-2	SELENIUM (DISSOLVED)	ug/l	1	U		U	2	U		U	2	U	2	U		U	9.3	J	1	U	6.8		7.8	ᆜ
7440-22-4	SILVER (DISSOLVED)	ug/l		U		U		U		Ų	1	U	1	U		U	1	U		U		U	1	U
7440-62-2	VANADIUM (DISSOLVED)	ug/l ug/l		U		U		U	• •	J		U	<u>1</u>	U		U	1 2.5	U	_	U J	1 2.7	U		J
7440-66-6	ZINC (DISSOLVED)	ug/l	2.1	J	5 l	U	5	U	1.6	J	5	U	5	U	1.2	J	2.5	U	2.4	J	2.1		2.3	J
75-25-2	BROMOFORM	ug/l	NA		NA	T	NA	П	NA	П	NA		NA		NA	Т	NA		NA		NA		NA	\neg
108-88-3	TOLUENE	ug/l			NA		NA		NA	寸	NA		NA		NA		NA		NA		NA		NA	\neg

Notes:
ug/L - micrograms per liter
U - not detected, laboratory reporting limit listed
J - concentration (or reporting limit) is estimated
R - data is rejected
DDA - Demolition Debris Area
NA - Not analyzed

Table 6-1 **Groundwater Quality** Phase II CSA Addendum **Demolition Debris Area** Bird Machine Company - Walpole, MA

															Down	gradie	nt of DDA											
		Station:			DD-MW-203				D	D-MW	-206				D	D-MW	<i>I-</i> 207							DD-MW-208				
	Field Sa	mple ID:	DD-MW-203-F	R04-X	DD-MW-203-R05	5-X	DD-MW-203-R06	3-X-E	DD-MW-206-R04	1-X [DD-MW-206-R0)5-X	DD-MW-207-R0-	4-X	DD-MW-207-R0	5-X	DD-MW-207-R0)5-D	DD-MW-207-R06-X	D-MW-208-R01-001	1- D Γ	D-MW-208-R01-0	01-X	DD-MW-208-R02-X	DD-I	MW-208-R03-X	DD-M	/W-208-R04-X
	Sam	ple Date:	6/5/2007	,	12/11/2007		5/19/2008		6/5/2007		5/19/2008		6/6/2007		12/11/2007		12/11/2007		5/19/2008	6/25/2007	T	6/25/2007		7/23/2007		12/11/2007	5	5/19/2008
		Depth:	4.5 - 12.5 fe	eet	4.5 - 12.5 feet		4.5 - 12.5 feet		5 - 11 feet		5 - 11 feet		5 - 15 feet		5 - 15 feet		5 - 15 feet		5 - 15 feet	3 - 13 feet	T	3 - 13 feet		3 - 13 feet		3 - 13 feet	3	3 - 13 feet
CAS Number	Analyte	Units	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result Q	Result Q	Į_	Result	Q	Result Q		Result Q	R	Result Q
51-28-5	2.4-DINITROPHENOL	/	5.0	UJ	5.1	UJ	5.1		F 2	111	5.1	U.J	5.1		5.1	UJ	5.5	U	5.2 U	5.1 R	+	5.1	R	5.2 UJ		5.1 U.		5.1 U
56-55-3	BENZO(A)ANTHRACENE	ug/l ug/l	5.2 0.31	U	_	U		IJ		UJ	0.3	U	-	UJ J	0.3	U	0.33	IJ	5.2 U 0.31 U		<u> </u>		U	5.2 UJ 0.31 U	_	0.3 U	+	0.3 U
50-32-8	BENZO(A)PYRENE	ug/i ug/l		U		U		U		U	0.2	U		U	0.3	U	0.33	IJ	0.31 U	0.0.	_		U	0.31 U	_	0.3 U	_	0.3 U
205-99-2	BENZO(B)FLUORANTHENE	ug/i	0.21	U		U		U		U	0.2	U	-	U	0.2	U	0.22	II	0.21 U		-		U	0.21 U	_	0.2 U		0.2 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l		U		U	***	U		U	0.51	U		U	0.51	Ü	0.55	U	0.51 U		_		Ü	0.21 J	_	0.51 U	_	0.51 U
207-08-9	BENZO(K)FLUORANTHENE	ug/I		U		U		U		U	0.3	U		U	0.3	Ü	0.33	U	0.32 U		_		Ü	0.21 J		0.3 U	_	0.3 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	- 3		U		U	***	J	5.2	U	0.49	J		U	5.1	Ü	5.5	U	0.51 J	0.0.			Ü	5.2 U	_	5.1 U		1.6 J
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.2	IJ	5.1	Ü	5.1	U	5.2	U	5.1	Ü	5.1	IJ	5.1	Ü	5.5	U	5.2 U		十		Ü	0.34 J		5.1 U	_	5.1 U
218-01-9	CHRYSENE	ug/l	1	II	• • • • • • • • • • • • • • • • • • • •	U		U	1	U	1	Ü	0.24	J	1	U	1.1	U	1 11	1 U	十		U	1 U		1 U	_	1 U
53-70-3	DIBENZ(A.H)ANTHRACENE	ug/l	0.52	U		U		U	0.52	U	0.51	Ü		U	0.51	Ü	0.55	U	0.52 U				U	0.42 J	_	0.51 U		0.51 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	5.2	U		Ü		U		U	5.1	Ü		U	5.1	U	5.5	Ü	5.2 U			0.0.	Ü	5.2 U		5.1 U	_	1.5 J
117-84-0	DI-N-OCTYL PHTHALATE	ug/l		Ü		Ü	0	Ü		Ü	5.1	Ü	_	J	5.1	Ü	5.5	Ü	5.2 U	0		• • • • • • • • • • • • • • • • • • • •	Ü	5.2 U		5.1 U		5.1 U
206-44-0	FLUORANTHENE	ug/l	1	Ü		Ü		U		U	1	Ü		J	1	Ü	1.1	Ü	1 U		亣	0.25	J	1 U		1 U	_	1 U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.52	Ü	0.51	Ū	0.51	Ū	0.52	U	0.51	Ū	0.51	Ü	0.51	Ū	0.55	U	0.52 U	0.51 U	٦	0.51	U	0.38 J		0.51 U		0.51 U
85-01-8	PHENANTHRENE	ug/l	0.21	Ü		Ū	0.2	Ū		U	0.2	Ū		Ū	0.2	Ū	0.22	U	0.21 U		٦	0.2	U	0.21 U		0.2 U	_	0.2 U
129-00-0	PYRENE	ug/l	5.2	Ū	5.1	Ū	5.1	Ū	5.2	U	5.1	Ū	0.29	J	5.1	Ū	5.5	U	5.2 U	0.28 J	ı T	5.1	U	5.2 U		5.1 U	_	5.1 U
		ŭ																			#							
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3	U	NA		NA	_	0.31	U	NA		0.02	U	NA		NA	\bot	NA	0.3 U	_	0.31	U	0.31 U		NA		NA
50-32-8	BENZO(A)PYRENE	ug/l		U	NA NA		NA NA		0.2.	U	NA NA	┢	0.2.	U	NA NA	┢	NA NA	+	NA NA	0.2 U	_	0.2.	U	0.21 U		NA	_	NA
205-99-2 191-24-2	BENZO(B)FLUORANTHENE BENZO(G.H.I)PERYLENE	ug/l ug/l	0.3	II	NA NA		NA NA	-	0.31 0.41	U	NA NA	1 1	*.*-	U	NA NA		NA NA	+	NA NA	0.3 U	_	0.0.	U II	0.31 U		NA NA	+	NA NA
207-08-9	BENZO(K)FLUORANTHENE	ug/I	0.4	Ü	NA NA		NA NA		Ŭ. I.	U	NA NA		**	IJ	NA NA		NA NA	1 1	NA NA	0.4 U		0	U	0.42 U		NA NA	+	NA NA
EPH1122	C11-C22 AROMATICS, ADJUSTED		100	Ü		U	NA		***	Ü	NA		****	Ü	100	U	100	U	NA	290 J	j	460	J	100 U		NA	-	100 U
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	100	U	NA		100	U	NA		110	U	100	U	100	U	NA	100 U	J	130		100 U		100 U		100 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.4	U	NA		NA		0.11	U	NA		U. 12	U	NA		NA		NA	0.4 U	J	0.11	U	0.42 U	_	NA	_	NA
206-44-0	FLUORANTHENE	ug/l	1	U	NA		NA			U	NA			U	NA		NA		NA	3.4	4	3.5		1 U	_	NA	-	NA
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.4	U	NA		NA		0.41	U	NA NA	1	VI.12	U	NA		NA NA	1	NA NA	0.4 U	#	0111	U	0.42 U		NA	-	NA
85-01-8	PHENANTHRENE	ug/l	0.2	U	NA		NA		0.21	U	NA		0.21	U	NA		NA		NA	0.62 J	4	0.37	J	0.21 U		NA		NA
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	11	U	1	U	1	UJ	1	U	1	U	1	U	1	U	1	U	1 U	1 U	J	1	U	NA		1 U		1 U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	1	U		U	1	U	1	U	1	U	5.7		10		10		3.9	1.7 J	工	1.7	J	NA		1 U		1 U
7440-39-3	BARIUM (DISSOLVED)	ug/l	12		11	_	11		26		16		38		45	$oxed{oxed}$	44	$oldsymbol{\perp}$	31	27	4	26		NA		25	_	31
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	2	U	1	U	1	U	1	U	1	U	2	U	2	U	1 U	2 U	4	1	U	NA		2 U	1	1 U
7440-43-9 7440-47-3	CADMIUM (DISSOLVED) CHROMIUM (DISSOLVED)	ug/l ug/l	1 16	U	1 15	U	1 0.63	U	1 23	U	<u>1</u> 0.51	U	<u>1</u> 4.9	U	1 10	U	<u>1</u> 91	U	1 U	1 U	4	1	U	NA NA	-	1 U 25	 	1 U 0.36 J
7440-47-3 7440-50-8	COPPER (DISSOLVED)	ug/i ug/l			NA	-+	0.63 NA	J	23 NA		0.51 NA	J	4.9 NA		NA	\vdash	9.1 NA	+	NA	9.8 NA	+	NA	-	NA NA		NA	_	0.36 J NA
7439-92-1	LEAD (DISSOLVED)	ug/I		U		U	1	U	1	U	1	U	1	IJ	1	U	1	IJ	1 U		十	1	U	NA NA	1	1 11	-	0.46 J
7439-97-6	MERCURY (DISSOLVED)	ug/l	•	U		U	0.2	Ü	0.2	U	0.2	Ü	0.2	Ü	0.2	Ü	0.2	U	0.2 U		十	0.2	Ü	NA NA	1	0.2 U	_	0.40 U
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.1			Ü	1.2		0.91	J	0.69	Ĵ		Ŭ	2	Ü	2	Ü	0.43 J		I	3.2		NA		0.96 J		2
7782-49-2	SELENIUM (DISSOLVED)	ug/l	1	U		U	1	U	1	U	1	U	1	U	1	U	1	U	1 U		J	2	U	NA		1 U		1 U
7440-22-4	SILVER (DISSOLVED)	ug/l	1	U		U		U		U	1	U	•	U	1	U	1	U	1 U		_		U	NA		1 U		1 U
7440-62-2	VANADIUM (DISSOLVED)	ug/l	1	Ų		Ų	0.20	J		U	1	Ų		U	2	Ų	2	U	1 U	0.0 .	4	1	J	NA		2 U	-	0.52 J
7440-66-6	ZINC (DISSOLVED)	ug/l	2.4	J	1.9	J	5.3		2.7		1.9	J	2.5	U	1.5	J	1.2	J	1.9 J	9.4	+	7.7	_	NA		20		110
75-25-2	BROMOFORM	ug/l	NA		NA		NA		NA		NA		NA		NA		NA		NA	5 U	ī	5	U	NA		NA		NA
108-88-3	TOLUENE	ug/l			NA		NA		NA		NA		NA		NA		NA		NA	5 U		5	U	NA		NA	_	NA

Notes:
ug/L - micrograms per liter
U - not detected, laboratory reporting limit listed
J - concentration (or reporting limit) is estimated

R - data is rejected
DDA - Demolition Debris Area
NA - Not analyzed

Table 11-1 Samples and Corresponding Sample Delivery Groups Phase II CSA Addendum Demolition Debris Area Bird Machine Company - Walpole, MA

Well	Field Sample ID	Sample Date	SDG
DD MW 004	DD-MW-001-R02-X	6/5/2007	360-10338
DD-MW-001	DD-MW-001-R02-X	6/5/2007	360-10338
	DD-MW-002-R01-X	6/5/2007	360-10381
	DD-MW-002-R01-X	6/5/2007	360-10381
DD MAY 000	DD-MW-002-R01-X	6/5/2007	360-10381
DD-MW-002	DD-MW-002-R01-X	6/5/2007	360-10381
	DD-MW-002-R02-X	5/19/2008	360-10338
	DD-MW-002-R02-X	5/19/2008	360-10338
	DD-MW-201-R04-X	6/5/2007	360-10381
DD MW 004	DD-MW-201-R04-X	6/5/2007	360-10381
DD-MW-201	DD-MW-201-R05-X	12/11/2007	360-10381
	DD-MW-201-R05-X	12/11/2007	360-10381
	DD-MW-201-R06-X	5/19/2008	360-10338
	DD-MW-201-R06-X	5/19/2008	360-10338
	DD-MW-201-R06-D	5/19/2008	360-10381
	DD-MW-201-R06-D	5/19/2008	360-10381
	DD-MW-203-R04-X	6/5/2007	360-10381
	DD-MW-203-R04-X	6/5/2007	360-10381
	DD-MW-203-R05-X	12/11/2007	360-10767
	DD-MW-203-R05-X	12/11/2007	360-10767
	DD-MW-203-R06-X	5/19/2008	360-10767
DD-MW-203	DD-MW-203-R06-X	5/19/2008	360-10767
	DD-MW-203-R06-X	5/19/2008	360-11279
	DD-MW-203-R06-X	5/19/2008	360-14013
	DD-MW-203-R06-X	5/19/2008	360-14013
	DD-MW-203-R06-X	5/19/2008	360-14013
	DD-MW-204-R02-X	6/5/2007	360-14013
DD 11111 004	DD-MW-204-R02-X	6/5/2007	360-14013
DD-MW-204	DD-MW-204-R05-X	5/19/2008	360-14013
	DD-MW-204-R05-X	5/19/2008	360-14013
	DD-MW-205-R03-X	6/6/2007	360-14013
DD MW 005	DD-MW-205-R03-X	6/6/2007	360-14013
DD-MW-205	DD-MW-205-R05-X	5/19/2008	360-14013
	DD-MW-205-R05-X	5/19/2008	360-14013
	DD-MW-206-R04-X	6/5/2007	360-14013
DD MW 000	DD-MW-206-R04-X	6/5/2007	360-14013
DD-MW-206	DD-MW-206-R05-X	5/19/2008	360-14013
	DD-MW-206-R05-X	5/19/2008	360-16608
	DD-MW-207-R04-X	6/6/2007	360-16608
	DD-MW-207-R04-X	6/6/2007	360-16608
	DD-MW-207-R05-X	12/11/2007	360-16608
DD 1111 007	DD-MW-207-R05-X	12/11/2007	360-16608
DD-MW-207	DD-MW-207-R05-D	12/11/2007	360-16608
	DD-MW-207-R05-D	12/11/2007	360-16608
	DD-MW-207-R06-X	5/19/2008	360-16608
	DD-MW-207-R06-X	5/19/2008	360-16608
	DD-MW-208-R01-001-X	6/25/2007	360-16608
	DD-MW-208-R01-001-X	6/25/2007	360-16608
	DD-MW-208-R01-001-D	6/25/2007	360-16608
	DD-MW-208-R01-001-D	6/25/2007	360-16608
DD-MW-208	DD-MW-208-R02-X	7/23/2007	360-16608
	DD-MW-208-R03-X	12/11/2007	360-16608
	DD-MW-208-R03-X	12/11/2007	360-16608
	DD-MW-208-R04-X	5/19/2008	360-16608
	DD-MW-208-R04-X	5/19/2008	360-16608
		=	

Notes:

SDG - Sample delivery group

Appendix A – Laboratory Analytical Reports (provided on compact disc)





DRAFT Method 3 Human Health and Environmental Risk Characterization

Demolition Debris Area, RTN 4-3024222 Former Bird Machine Company Walpole, MA

Submitted to:

Baker Hughes Incorporated Houston, TX

Submitted by:

AMEC Earth & Environmental, Inc. Westford, Massachusetts

November 2011



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LIST OF ACRONYMS

ADD Average Daily Dose
ADE Average Daily Exposure

AIS Asbestos in Soil

AUL Activity and Use Limitation
bgs Below Ground Surface
BMC Bird Machine Company
COC Chemical of Concern

COPC Chemical of Potential Concern

COPEC Chemical of Potential Environmental Concern

CSA Comprehensive Site Assessment

CSF Cancer Slope Factor
CSM Conceptual Site Model

d Day

DDA Demolition Debris Area

DWF Division of Fisheries and Wildlife EcoSSL Ecological Soil Screening Level

EH 855 Estimated Habitat 855

ELCR Excess Lifetime Cancer Risk EPC Exposure Point Concentration

EPH Extractable Petroleum Hydrocarbon ERC Environmental Risk Characterization

ft Feet

GI Gastrointestinal GW Groundwater

HHRC Human Health Risk Characterization

HI Hazard Index HQ Hazard Quotient

hr Hour

IRIS Integrated Risk Information System

kg Kilogram

LADD Lifetime Average Daily Dose

LOAEL Lowest Observed Adverse Effect Level

LRA3 Lead Release Area 3

MADEP Massachusetts Department of Environmental Protection

MBA Manufacturing Building Area
MCP Massachusetts Contingency Plan

MESA Massachusetts Endangered Species Act

MFG Million Fibers per Gram

mg Milligram

MMCL Massachusetts Maximum Contaminant Level

m³ Cubic Meter µg Microgram

NHESP Natural Heritage and Endangered Species Program

NOAEL No Observed Adverse Effect Level

NSR No Significant Risk

PAH Polycyclic Aromatic Hydrocarbon



LIST OF ACRONYMS, continued

PCB Polychlorinated Biphenyls
PH 1072 Priority Habitat 1072
PM Particulate Mass
PM10 Respirable fraction

RAF Relative Absorption Factor
RC Risk Characterization
RfC Reference Concentration

RfD Reference Dose

RTN Release Tracking Number

SF Scaling Factor SRS South Rail Spur

SVOC Semi-volatile Organic Compound TEF Toxicity Equivalency Factor

TEQ Toxic Equivalent

TRV Toxicity Reference Value UCL Upper Concentration Limit

URF Unit Risk Factor

USEPA United States Department of Environmental Protection

UTL Upper Tolerance Limit
VOC Volatile Organic Compound
VPH Volatile Petroleum Hydrocarbon

WOE Weight-of-Evidence WPA Wetlands Protection Act



1.0 INTRODUCTION

The Bird Machine Company (BMC) facility property ("the BMC facility" or "the facility") occupies approximately 134 acres of land located in Walpole, Massachusetts. The BMC facility, which formerly manufactured machinery, has been assigned multiple Release Tracking Numbers (RTNs) under the Massachusetts Contingency Plan (MCP). This risk assessment addresses the release at the Demolition Debris Area (DDA), which was assigned RTN 4-3024105 and was classified as a Tier II Disposal Site in July 2005 (Weston, 2007). The DDA RTN was linked to RTN 4-3024222 in the January 2008 Tier 1B Permit Application for the facility. Bird Machine Company is no longer in operation, and most of the buildings have been permanently removed. The location of the property is depicted in **Figure 1-1**.

A draft Phase II Comprehensive Site Assessment (Phase II CSA) Addendum has been completed for the DDA. The Phase II CSA Addendum addresses volatile organic constituents (VOCs), semi-volatile organic constituents (SVOCs), volatile petroleum hydrocarbons (VPH), extractable petroleum hydrocarbons (EPH), polycyclic aromatic hydrocarbons (PAHs), dioxin/furan congeners, and various metals detected in soil and groundwater samples collected from the DDA. The Phase II CSA Addendum also includes evaluations of asbestos in soil (AIS) identified at the DDA.

This Risk Characterization (RC) has been prepared by AMEC Earth & Environmental, Inc. of Westford, Massachusetts (AMEC) to support the Phase II CSA Addendum. The previous reports prepared for the DDA by Weston Solutions, Inc. (Weston) of Concord, New Hampshire, contain information on regulatory compliance, the methods and findings of site assessment activities, and preliminary and comprehensive response actions conducted at the DDA. Information in these reports was used to prepare this risk characterization. Data from site investigations completed by AMEC, site assessment activities completed by Weston, and information from other sources (e.g., Massachusetts Department of Environmental Protection [MADEP] and United States Environmental Protection Agency [U.S. EPA]), were used to complete the RC.

In accordance with the requirements of 310 CMR 40.0000 Subpart I of the MCP, the Method 3 RC addresses risk of harm to human health, public welfare, safety, and the environment. This RC has been conducted assuming that an Activity and Use Limitation (AUL) will be implemented at the DDA prohibiting future development at the site and the disturbance of surface soil. Therefore the RC does not evaluate DDA use or development other than incidental trespassing.

The RC conforms with the requirements of 310 CMR 40.0000 Subpart I of the MCP, and the Massachusetts Department of Environmental Protection's (MADEP's) *Guidance for Disposal Site Risk Characterization - In Support of the Massachusetts Contingency Plan* (MADEP, 1995),



to evaluate potential risk of harm to human health, safety, public welfare and the environment posed by DDA conditions.

A Method 3 approach has been used to assess both human health and environmental risk. The Method 3 human HHRC has four steps. The first step, Hazard Identification, involves identification of the constituents of potential concern (COPCs) detected at the DDA. The second step, Dose response Assessment, describes the relationship between the magnitude of exposure for each COPC (dose) and the occurrence of health effects (response). The third step, Exposure Assessment, identifies of potential human receptors based on characteristics of the DDA and the surrounding area. Subsequently, the magnitude and frequency of receptors' potential exposure to COPCs is quantified. The fourth step, Risk Characterization, combines the information from the Exposure Assessment with the information from the Dose response Assessment to determine the likelihood of adverse non-carcinogenic health effects or excess lifetime carcinogenic effects for each receptor for each potential exposure pathway identified in the Exposure Assessment. The risks associated with each exposure pathway are summed to obtain an estimate of total risk for each receptor. Details on these steps of the HHRC, as well a characterization of risk to safety and public welfare, are provided in **Section 3.0** of this report.

Environmental risk characterizations (ERCs) typically consist of two phases: Stage I environmental screening and Stage II risk characterization. Environmental screening is designed to determine whether a more detailed evaluation is necessary based on three criteria: 1) whether environmental receptors could potentially be exposed to constituents at the DDA presently or in the future; 2) whether significant environmental harm is "readily apparent" for each of the potential exposures identified; and 3) whether any of the potential exposure pathways could result in "potentially significant" exposures. Since significant current or potential future exposures to constituents in DDA media were not ruled out by the environmental screening, a Stage II ERC was performed. The analysis was conducted in accordance with the MADEP's *Guidance for Disposal Site Risk Characterization* (MADEP, 1995; 1996).

Details of the risk characterization are presented in the remainder of this report, as follows: **Section 2.0** provides a brief overview of the DDA setting and history. **Section 3.0** presents the risk characterization of human health, safety, and public welfare. The environmental RC is presented in **Section 4.0**. The summary and conclusions are presented in **Section 5.0**. A list of reference materials used to conduct the RC follows **Section 6.0**.



2.0 SITE BACKGROUND

The BMC facility manufactured machinery for the paper industry and various other industries from 1920 until 2004, when the company ceased manufacturing operations at the property. Most of the buildings and equipment were then removed from the property, although some materials (such as pieces of concrete) remain.

For the purposes of evaluating potential exposures and risks, the BMC was divided into several exposure areas: the South Rail Spur (SRS) is in the southwestern section of the property; the lead Release Area 3 (LRA3) is in the eastern section of the property (which encompasses a fill area); the Manufacturing Building Area (MBA), which also contains Lead Release Area 1 and Lead Release Area 2, in the southeastern portion of the property (where most of the structures used during operations at the facility were housed); the Neponset River, which runs along the eastern boundary of the property; and the DDA (as previously defined, the Demolition Debris Area) (the "Site" as defined in this report) in the northwestern portion of the BMC property.

Figure 2-1 shows the DDA in relation to the other areas of the BMC. The DDA consists of three contiguous clearings known as the Eastern, Western and Central Clearing areas, which served as a disposal area dating back possibly to the 19th century. Material in the DDA is primarily inorganic fill, demolition debris, and manufacturing/product testing wastes. Since the most recent removal action (see below), the area has been undisturbed.

Various Release Abatement Measures and Immediate Response Actions have been performed throughout the BMC. The Phase I and Phase II reports for these various areas detail the regulatory and remedial history. The DDA was classified as a Tier II Disposal Site in July 2005. In September 2005, Weston observed fibrous material suspected of containing asbestos. Over the next few months, Weston and its subcontractors confirmed asbestos and conducted a series of removals as an Immediate Response Action. Weston concluded that asbestos could be visibly identified. Following the final removal, the excavation area was lined with geotextile and covered with soil.

A Phase II CSA was prepared by Weston in July 2007. The Phase II CSA included a Method 1 RC. The Phase II CSA was unable to conclude No Significant Risk (NSR) due to, among other things, the visually observed presence of asbestos in soil. A Phase III RAP was prepared by Weston in July 2007 which selected a soil cover remedy for implementation. Environmental risks were not characterized at the time. Shortly after these reports were finalized, the DDA was combined with other RTNs in the Tier IB permit.

Additional sampling was conducted in 2011 to delineate the nature and extent of the asbestos contamination at the DDA. With this additional data, this Method 3 RC was prepared to replace the previous Method 1 RC, and to provide environmental risk characterization for the DDA.



3.0 HUMAN HEALTH RISK CHARACTERIZATION

This section presents the methodology used to evaluate potential risks to human health using a Method 3 approach. Section 3.1 presents the Hazard Identification and Section 3.2 presents the Dose Response Assessment. The Exposure Assessment is described in Section 3.3; and the Risk Characterization is presented in Section 3.4. Uncertainties associated with the human health risk characterization are discussed in Section 3.5.

3.1 Hazard Identification

This section includes a description of the data used as the basis of the risk characterization and presents the selection of contaminants of potential concern (COCs) in each medium.

3.1.1 Database

The DDA has previously been determined to be limited to the upland area shown on Figure 2-1 and does not extend to Cedar Swamp Brook (Weston, 2007). The database for this RC consists of soil, groundwater, and tissue (earthworm) samples as follows:

- <u>Soil</u> Because an AUL prohibiting intrusive activities will be instituted, only surficial soil
 [0-3 feet (ft) below ground surface(bgs)] was considered. Soil data that were included in
 the DDA RC exposure point concentrations (EPCs) consist of results obtained from
 sampling events performed in December 2004; May, November, and December 2005;
 June, September and October 2006; and May 2007.
- Asbestos in soil An asbestos in soil (AIS) delineation program was performed for surficial soil in April 2011. In June 2011, three samples from the one sample in which asbestos was identified underwent elutriator analysis to estimate an airborne fiber concentration.
- <u>Groundwater</u> Multiple sampling events for groundwater have been performed dating back to 2006. The most recent data set (collected in May 2008) is used in this RC.
- Tissue (earthworm) collected September 2006.

Data have been reviewed and are of suitable quality for inclusion in site characterization. A representativeness and data usability assessment are presented in Section 11.0 of the Phase II.

For both soil and groundwater data, non-detect results for constituents detected in at least one sample in a medium and exposure area were assumed to equal one-half of the analytical reporting limit. Field duplicates were averaged using the following methodology:



- 1. When both samples in the pair of duplicates had detected values, the results from the primary and its duplicate were averaged and treated as one detected concentration at the location:
- 2. When only one of the duplicate samples had a detected value, the detected value was averaged with half the reporting limit of the non-detected value and the average was treated as one detected concentration at the location; and
- 3. When both samples had non-detect values, half of the lower of the two reporting limits was used as the non-detected concentration at the location.

When analytes were analyzed using multiple methods (for example, naphthalene from EPH and SVOC analyses), one result was selected using the following methodology:

- 1. When both samples in the pair of duplicates had detected values, the higher result from was selected and treated as one detected concentration at the location;
- 2. When only one of the duplicate samples had a detected value, the detected value was selected and treated as one detected concentration at the location; and
- 3. When both samples had non-detect values, the lower reporting limit was selected and treated as one non-detect concentration at the location.

The datasets used in this risk characterization are presented in **Attachment A**. Soil, groundwater, and earthworm sample locations are presented on **Figure 2-2**. AlS sampling locations are presented on **Figure 2-3**.

In total 78 soil samples, including duplicates, were collected from 57 locations in DDA between 12/20/2004 and 06/6/2007. Only those soil samples from the 0 to 3 ft bgs depth interval were used in the RC. Groundwater samples from eight shallow wells were collected on 05/19/2008. All data collected from these wells were used in the RC.

3.1.2 Selection of Contaminants of Concern

Soil: Any constituents detected at least once at a concentration exceeding background in an exposure area were included as a COC in that medium and exposure area. For metals and PAHs with background concentrations published in the MassDEP's (MADEP, 2002a) Technical Update entitled "Background Concentrations of Polycyclic Aromatic Hydrocarbons and Metals in Soil," the concentration in "natural" soil was considered background. For all other constituents, any concentration detected above the laboratory reporting limit was considered above background.

Groundwater: Because there are no published background concentrations for groundwater, any constituents detected in the most recent sampling round for groundwater were considered a COC. However, a background evaluation was completed for arsenic, which is ubiquitous and was detected at a relatively low concentration (4 ug/l) in one downgradient well in the most



recent sampling round. The background evaluation was based on groundwater data collected in other areas of the BMC determined to be free of influence from the DDA. ProUCL was used to calculate an upper tolerance limit (UTL), which represents the upper end of a fixed proportion of the population with a stated confidence, in this case the 90th percentile with 95% confidence. In other words, the UTL is the value above which, with 95% certainly, only 10% of the values in the true population fall. The UTL calculated by ProUCL using the background arsenic data for wells at the facility was 14.3 ug/L. This evaluation appears in **Attachment B.** Based on the background analysis, arsenic was excluded as a COC as it was detected below background concentrations. All other constituents detected in the 2008 sampling round at concentrations above the laboratory reporting limits were selected as COCs for groundwater, with the exception of 2,4-dinitrophenol, di-n-butylphthalate, and lead, which were eliminated because of low frequency of detection and low concentration.

Tables 3-1 and **3-2** present the selection of COCs for DDA soil and groundwater, respectively. Asbestos is also a COC; the estimated asbestos fiber concentrations is described in Section 3.3.8.3.

3.2 Dose Response Assessment

The purpose of the Dose Response Assessment is to identify the relationship between the quantity of COCs to which receptors may be exposed (dose) and the likelihood of an adverse health effect (response). Both noncarcinogenic (i.e., threshold) and carcinogenic (i.e., non-threshold) health effects were considered in the dose response assessment. The information provided in the Dose Response Assessment was combined with the results of the Exposure Assessment to provide an estimate of potential health risk. Noncarcinogenic dose response information is presented in Section 3.2.1, and Section 3.2.2 discusses carcinogenic dose response.

Dose response information used in this RC was obtained from MADEP and EPA publications. Toxicity values for EPH and VPH carbon fractions were obtained from the MADEP Policy #WSC-02-411, *Implementation of the MADEP VPH/EPH Approach* (MADEP, 2002b). Toxicity values for other COCs were obtained from MADEP's MCP Toxicity.xls spreadsheet used by MADEP to derive Method 1 Standards (MADEP, 2009). References in the MADEP spreadsheets to U.S. EPA data, including the Integrated Risk Information System (IRIS), were checked and updated, as necessary. The toxicity data used in the Method 3 RC of soil are shown in **Table 3-3**.

3.2.1 Noncarcinogenic Dose Response

Constituents with known or potential noncarcinogenic effects are assumed to have a dose below which no adverse effect occurs, or conversely, above which an effect may be seen. In laboratory experiments, this dose is known as the "No Observed Adverse Effect Level"



(NOAEL). The lowest dose at which an adverse effect is seen is called the "Lowest Observed Adverse Effect Level" (LOAEL). By applying uncertainty factors to the NOAEL or the LOAEL, Reference Doses (RfDs) or Reference Concentrations (RfCs; for air) are developed for chronic and, in some cases, subchronic exposures to constituents with potential noncarcinogenic effects. Many of the non-carcinogenic dose response values provided by MADEP (2009) were developed by the U.S. EPA and are reported in U.S. EPA (2010a,b), while other values provided in MADEP (2009) were developed or selected by MADEP.

Uncertainty factors account for uncertainties associated with the dose response data, such as the appropriateness of using an animal study to derive a human dose response value, and the potential for especially sensitive subpopulations to exist, which may not be adequately represented by the laboratory test animals. For constituents with potential noncarcinogenic effects, the RfD/RfC provides reasonable certainty that, if the specified exposure dose is below the threshold, no noncarcinogenic health effects are expected to occur. RfDs are expressed in terms of milligrams of constituent per kilogram of body weight per day (mg/kg-day) and are used to evaluate estimated oral and dermal exposures. RfCs are expressed as milligrams per cubic meter (mg/m³) and are used for inhalation. RfDs and RfCs are sometimes inter-converted. **Table 3-3** summarizes the toxicity values for the COCs evaluated here by the ingestion, dermal, and inhalation exposure routes.

3.2.2 Carcinogenic Dose Response

The U.S. EPA assumes for regulatory risk assessment that no threshold dose exists (U.S. EPA, 1997b, 2010a). In other words, U.S. EPA assumes that a finite level of risk may be associated with any dose above zero. In March 2005, U.S. EPA issued new cancer guidelines (U.S. EPA, 2005), the purpose of which is to recommend principles and procedures to guide U.S. EPA scientists in assessing the cancer risks from constituents or other agents in the environment when deriving toxicity values. U. S. EPA uses a two-part system for characterizing the extent to which the available data support the hypothesis that an agent causes cancer in humans.

U.S. EPA's first step in evaluating a potential carcinogen is to assign a weight-of-evidence (WOE) classification. Under U.S. EPA's previous cancer guidelines released in 1986, the WOE was described by categories "Group A" through "Group E," with Group A category reserved for known human carcinogens, while Group E category was the other end of the spectrum, representing constituents/agents with evidence of non-carcinogenicity. In the U.S. EPA's more recent approach for carcinogen risk assessment (U.S. EPA, 2005), all scientific information is considered in determining whether and under what conditions an agent may cause cancer in humans. Furthermore, the WOE provides a narrative approach to characterize carcinogenicity rather than distinct categories by summarizing the evidence about the likelihood of the constituent being a human carcinogen. Five standard WOE descriptors are currently used as part of the narrative, including:



- 1. Carcinogenic to Humans;
- 2. Likely to be Carcinogenic to Humans;
- 3. Suggestive Evidence of Carcinogenic Potential;
- 4. Inadequate Information to Assess Carcinogenic Potential; and
- 5. Not Likely to be Carcinogenic to Humans.

As part of the updated guidance on evaluating potentially carcinogenic constituents, the U.S. EPA emphasizes the value of understanding the biological changes that the agent of interest can cause (e.g., mode of action) and how these changes might lead to the development of cancer. This information, as well as the agent's human carcinogenic potential, is to be described in a narrative prepared by U.S. EPA's scientists, summarizing the full range of available evidence and describes any conditions associated with conclusions about an agent's hazard potential, including which populations or life stages may be particularly susceptible. Since the data for many of the potentially carcinogenic constituents have not been re-evaluated since the initial derivation of the cancer slope factors under the 1986 cancer guidelines, the cancer toxicity information presented in Integrated Risk Information System (IRIS - U.S. EPA's database of recommended cancer slope factors and reference doses for use in risk assessments) represents cancer toxicity information derived under the 1986 guidelines, with more recent cancer evaluations conducted for a limited number of constituents under the more recent 2005 guidance (U.S. EPA, 2010a).

The second step in the carcinogenicity evaluation process is the calculation of a quantitative estimate of carcinogenic potency. The U.S. EPA has developed computer models that extrapolate the observed responses at high doses used in animal studies to predict responses in humans at the low doses encountered during environmental exposures. The models developed by the U.S. EPA assume no threshold and usually consider animal (and sometimes human) data to estimate carcinogenic potency. Further, the models assume that carcinogenic dose response is linear at low doses. U.S. EPA refers to this numerical estimate of the dose response factor (or the slope of the line plotted from dose vs. response) as the cancer slope factor (CSF) for oral exposures. A CSF is expressed in terms of the inverse of milligrams of agent per kilogram body weight per day [(mg/kg-day) $^{-1}$] and represents the upper-bound excess lifetime cancer risk estimate that results from a daily exposure to an agent at a dose of 1 mg/kg-day. A Unit Risk Factor (URF) represents the Excess Lifetime Carcinogenic Risk (ELCR) per microgram per cubic meter (μ g/m 3) of contaminant in air and is expressed as (μ g/m 3) $^{-1}$. CSFs and URFs are sometimes inter-converted.

Table 3-3 summarizes carcinogenic toxicity values for COPCs used in this risk characterization.

3.3 Exposure Assessment

This section identifies the type and magnitude of potential exposures to COPCs that may occur at the DDA under current and reasonably foreseeable future use. First, potential receptors are



identified based on conditions present at the DDA and surrounding area. Next, potential routes and pathways of exposure are identified for each receptor, based on information about activities that typically occur or may occur in the area. Following these steps, EPCs are estimated, and potential exposures are quantified for receptors.

3.3.1 Site Use and Activities

The DDA is located within the larger BMC property, which is inactive. Manufacturing operations at the property were discontinued in 2004, and most buildings associated with the property have been demolished. Fencing is present at a portion of the property. A security guard is on duty (at the entrance of the property) to prevent unauthorized vehicle access and to report trespassing activities to the owners and municipal authorities. No development of the DDA is currently planned.

The DDA was a disposal area at the northwest part of the property and remains undisturbed. Due to the fencing, current receptors at the facility, if any, are limited to occasional trespassers. This RC assumes an Activity and Use Limitation (AUL) will be implemented at the DDA to prevent disturbance. As such, future site use is considered to be limited to occasional trespassers as well.

3.3.2 Identification of Receptors

Consistent with the requirements of the MCP, 310 CMR 40.0923, the exposure assessment considered both current and reasonably foreseeable future Site activities and uses. The only potential current receptors are local resident adolescents who may trespass on the property and the security personnel. The existing fencing, waterways, and active railways surrounding the Site are expected to effectively prevent young children from gaining access to the Site.

This Method 3 RC assumes that an AUL will be implemented on the property that would prohibit disturbance of the soil at the DDA and excavation, grading or development for any purpose. As such, an occasional trespasser is the only current and future receptor.

3.3.3 Identification of Exposure Routes and Pathways

Exposure pathways are the mechanisms by which receptors may be exposed to COCs at the Site. According to MADEP (1995), the following elements must be present in order for a potential human exposure pathway to be complete:

- 1. a constituent source:
- 2. a mechanism by which a constituent may be released to the environment;
- 3. an environmental transport medium;
- 4. an exposure point (discussed above); and,
- 5. a receptor with a route of exposure at the point of contact (discussed above).



Under current and foreseeable future conditions, trespassers are the only receptors that may be exposed to COCs in DDA soil. The potentially complete exposure pathways for the trespasser include incidental ingestion, dermal contact, and inhalation of particles (dust); inhalation of asbestos fibers from DDA soil is also considered. The Town of Walpole has designated the BMC property as located within a groundwater Recharge Zone (BSWC, 2007). As such, although groundwater at the Site is not currently used nor is likely to be used in the future as a source of drinking water, use of Site groundwater as drinking water must be considered a potential future exposure pathway.

3.3.4 Soil and Groundwater Categorization

Based on current conditions, the most likely potential receptors are local resident adolescents who may trespass on the property. The waterways and active railways surrounding the Site are expected to effectively prevent young children from gaining access to the Site. The activity of adolescent trespassers is characterized as low frequency/low intensity, because access to the Site would similarly be limited by the fencing. The fencing only surrounds part of the Site, but it is still expected to have some effect as a deterrent when it comes to adolescent trespassers. Trespassers would not likely be engaged in activities that would result in contact with deeper soil. Although there are currently no plans for redevelopment of the property, this risk characterization was conducted assuming that an AUL will be implemented at the Site prohibiting future disturbance of surface soil. Based on this assumption, soil from 0 to 3 feet is categorized as S-2 and soil from 3 to 15 ft bgs is categorized as S-3.

As mentioned above, the Town of Walpole has designated the Site as a groundwater Recharge Zone (BSWC, 2007). As a result of this designation, the Town's Board of Sewer and Water Commissioners has indicated its belief that "the Site is in a Potential Drinking Water Source area" according to conditions specified at 310 CMR 40.0932(4). For this reason, groundwater in all exposure areas is categorized as GW-1. Because no buildings are present at the Site and future buildings will be prohibited by the AUL, groundwater category GW-2 is not applicable. Site groundwater is also categorized as GW-3 because, as described in 310 CMR 40.0932, this category is applicable to all groundwater in the Commonwealth, as all groundwater is assumed to potentially discharge to surface water.

3.3.5 Hot Spot Evaluation

The MCP requires identification and evaluation of "hot spots." As defined by the MCP in 310 CMR 40.0006, a hot spot is "a discrete area where the concentrations of oil or hazardous material or the thickness of Nonaqueous Phase Liquid are substantially higher than those present in the surrounding area. A hot spot shall be identified based on consideration of both the concentrations or thickness of an oil or hazardous material within a contaminated area and the spatial pattern of that contamination. The areal extent and spatial pattern of a hot spot may



be determined through the analytical results from multiple samples taken within the area, or the results of limited sampling in combination with other knowledge about the release, such as the presence of discoloration, odors or a defined source area." One method for evaluating whether an area is a hot spot includes comparison of OHM concentrations in an area with the concentrations in the surrounding area. An area that has a concentration greater than the applicable Method 1 standard and either 10 times (if the area may be contacted preferentially or more frequently) or 100 times greater than the concentration in the surrounding area. Since the Site is currently unoccupied and (although there are currently no plans for development of the property) future development could occur anywhere on the Site, and because no areas of the Site appear especially attractive to potential trespassers or would be expected to have higher exposure frequency than other areas, hot spots were defined as areas with concentrations 100 times higher than the concentration of the surrounding area. Hot spots may be a single sample, multiple samples from the same soil boring, or samples from multiple contiguous soil borings over a small area. Hot spots were identified using the methodology described in the following paragraphs.

Maximum detected concentrations in soil in each exposure area were first compared to Method 1 S-1/GW-1 soil standards. Any constituent with a maximum detected concentration in an exposure area below its Method 1 standard was excluded from further evaluation. Constituents with maximum concentrations above the S-1/GW-1 standard were further evaluated by plotting the highest detected concentrations on a map of the exposure area to determine if these high concentrations were scattered or clustered. Isolated high concentrations in a single sample were evaluated by comparing those concentrations to surrounding concentrations. When high concentrations were clustered, the concentrations were examined to determine if they extended over multiple contiguous sample locations. In such cases, the average concentrations of constituents in the samples with high concentrations were estimated and compared to the surrounding concentrations.

If concentrations were not more than 100 times higher than the concentrations in the surrounding area, they were not identified as "hot spots." Hot spot analysis of groundwater as a potential water supply was not performed because each well is considered a distinct exposure point. No hot spots were identified in DDA soil.

3.3.6 Identification of Exposure Points

As indicated in Section 3.1.1, data that are representative of current conditions in soil were used to develop EPCs in soil at the DDA. Soil from a depth of 0 to 3 ft bgs is considered to be accessible and therefore constitutes the current soil exposure zone. Due to the AUL that will be implemented for the Site, the only soil that will be available in the future is also surficial (0 to 3 feet).



3.3.7 Estimation of Exposure Point Concentrations

The methods used to calculate EPCs in soil and groundwater are described in the sections below. In the special case of dioxins/furans, one additional data management step is necessary to calculate EPCs. Toxicity values and Method 1 standards are not available for all of the individual dioxin/furan congeners included in the analytical method, but are available for 2,3,7,8-tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD). The toxicity of other dioxin/furan congeners has been estimated relative to the toxicity of 2,3,7,8-TCDD. In each sample analyzed for dioxins and furans, the concentration of 2,3,7,8-TCDD toxic equivalents (TEQ) was estimated by first multiplying the concentration of each congener by the toxicity equivalency factor (TEF) for that congener developed in 2005 by the World Health Organization (Van den Berg et al., 2006), then summing the toxic equivalent concentrations for the congeners. The 2005 WHO TEFs are presented in **Table 3-4**. Any congener not detected in a sample was assumed to be present at one half the detection limit in the computation of the TEQ concentration in that sample.

3.3.7.1 Soil EPCs

The arithmetic mean of COC concentrations was used to represent the EPC in each exposure area. Arithmetic mean concentrations were calculated assuming that non-detected constituents were present at one-half the sample quantitation limit. In accordance with the MCP, arithmetic mean concentrations are used as EPCs unless there is evidence of tremendous skew in the data, since it is unlikely that the arithmetic averages would underestimate the true means. It is our opinion that the sampling program conducted at the Site provides sufficient analytical data to estimate EPCs that are representative of Site conditions. Specifically, use of average concentrations is justified based on the following:

- A review of the data indicates that the exposure areas have been adequately characterized based on the number of sampling locations. Qualitative evaluation of sampling density indicates that based on the size of the exposure area, a sufficient number of samples have been collected to provide adequate areal coverage of the release area(s); and
- The sampling strategy employed at the exposure areas results in a conservative assessment of soil quality conditions as sampling locations are biased to areas of higher concentration (the release areas).

The EPCs for soil were calculated from samples collected for the 0 to 3 ft bgs depth, and are presented in **Table 3-1**. The locations of the soil samples are shown in **Figure 2-1**.

3.3.7.2 Groundwater EPCs

As required by the MCP, each groundwater well was considered a hypothetical potential source of future drinking water, and each was considered a separate exposure area. Therefore, all



groundwater wells sampled in 2008 were included, regardless of groundwater depth. EPCs in individual wells were estimated using the sample results for that well. EPCs for groundwater are presented in **Table 3-2**. The locations of the groundwater samples are shown in **Figure 2-1**.

3.3.8 Quantification of Potential Exposures

This section describes the equations and assumptions used to evaluate potential exposures to COPCs in the Method 3 evaluation at DDA. These equations are consistent with equations presented by MADEP (1995).

The Average Daily Dose (ADD) was calculated to estimate a receptor's potential daily intake from exposure to constituents with potential noncarcinogenic effects. According to MADEP (1995), the exposure dose should be calculated by averaging over the period of time for which the receptor is assumed to be exposed. The ADD for each constituent via each route of exposure was then compared to the noncarcinogenic toxicity value (that is, the RfD) for that constituent in order to estimate the potential noncarcinogenic hazard index due to exposure to that constituent via that route of exposure.

For constituents with potential carcinogenic effects, the lifetime average daily dose (LADD) was calculated to estimate potential exposures over the course of a lifetime (70 years). Subsequently, the LADD for each constituent via each route of exposure was multiplied by the CSF for that constituent to estimate the potential carcinogenic risk due to exposure to that constituent via that route of exposure.

The equations used to estimate ADDs and LADDs are presented below. The human exposure parameter values used in each potential exposure pathway are presented in **Table 3-6**, while **Table 3-5** summarizes certain constituent specific factors required to implement the equations (i.e., relative absorption factors). The spreadsheets used to calculate ADD and LADD from these equations and parameter values are contained in **Attachment D**.

3.3.8.1 Soil

Exposure to soil was assumed to occur via incidental ingestion, dermal contact, and inhalation of particles (fugitive dust). ADDs and LADDs for soil ingestion were calculated as follows:

ADD
$$_{ing}$$
 or LADD $_{ing}$ =
$$\frac{C_{soil} \times IR_{soil} \times FI \times RAF_{os} \times CF \times EF \times EP}{AP \times BW}$$

where:

ADD_{ina} = Average Daily Dose Due to Ingestion (mg/kg-day)

LADD_{ing} = Lifetime Average Daily Dose Due to Ingestion (mg/kg/day)



C_{soil} = Constituent Concentration in Soil (mg/kg)

 IR_{soil} = Soil Ingestion Rate (mg/day)

FI = Fraction of Soil Ingested from the Site (unitless)
RAF_{os} = Relative Absorption Factor (Oral-Soil) (unitless)

CF = Conversion Factor (10^{-6} kg/mg) EF = Exposure Frequency (days/year)

EP = Exposure Period (years)

BW = Body Weight (kg)

AP = Averaging Time (EP x 365 days/yr, ADD; 70yr x 365 days/yr, LADD)

ADDs and LADDs for dermal absorption were calculated as follows:

ADD _{der} or LADD _{der} =
$$\frac{C_{soil} \times SA \times AF \times RAF_{ds} \times CF \times EF \times EP}{AP \times BW}$$

where:

 ADD_{der} = Average Daily Dose Due to Dermal Contact (mg/kg-day)

LADD_{der} = Lifetime Average Daily Dose Due to Dermal Contact (mg/kg/day)

C_{soil} = Constituent Concentration in Soil (mg/kg) SA = Skin Surface Area Exposed (cm²/day) AF = Soil to Skin Adherence Factor (mg/cm²)

 RAF_{ds} = Relative Absorption Factor (Dermal-Soil) (unitless)

 $CF = Conversion Factor (10^{-6} kg/mg)$ EF = Exposure Frequency (days/year)

EP = Exposure Period (years)

BW = Body Weight (kg)

AP = Averaging Period (EP x 365 days/yr, ADD; 70yr x 365 days/yr, LADD)

3.3.8.2 Inhalation of Particulates

Exposure via inhalation of soil-derived fugitive dust is a function of the concentration at the source (e.g., soil), frequency and duration of contact, and a factor describing the concentration of respirable particles in air.

MADEP (2008) considers that potential exposure via inhalation of dust occurs via two uptake pathways: uptake by the gastrointestinal (GI) tract following coughing up and subsequent swallowing of particulates trapped by the mucosa of the upper respiratory track and uptake by the respiratory system following inhalation into the lungs. To calculate the exposure associated with these two uptake pathways, MADEP assumes the following for the construction worker scenario (which AMEC conservatively also used for the trespasser and utility worker):

- 100% of respirable particulate mass (PM) is equal to or less than 30 microns in diameter (<=PM₃₀)
- 40% of total respiratory particulate mass is equal to or less than 10 microns in diameter (<=PM₁₀)



- 100% of inhaled particulates greater than 10 microns but less than or equal to 30 microns are swallowed.
 50% of inhaled particulates equal to or less than 10 microns are swallowed
- 50% of inhaled particulates equal to or less than 10 microns enters the lungs.

Based on these assumptions, the effective exposure concentration of respirable particulates for the GI system is 1.5 times the concentration of PM_{10} , while that for the lungs is 0.5 times the concentration of PM_{10} . Using these effective exposure concentrations, compound average daily doses for the GI and respiratory systems can be estimated using the following equations.

Average Daily Dose for the GI System (ADD_{inhal-GI}):

$$ADD_{inhal} = \frac{C_{part} \times 1.5 \times PM_{10} \times IR_{air} \times RAF_{i} \times ET \times EF \times EP \times CF}{AP \times BW}$$

where:

ADD _{inhal-Gl} = Average Daily Dose due to coughing up and subsequent ingestion of inhaled particulates; expressed in mg/kg-day

[C_{part}] = Concentration of constituent in airborne particulates

 $[PM_{10}]$ = Concentration in air of particulates less than or equal to 10 microns in diameter

IR_{air} = Inhalation rate for the receptor of concern during the period of exposure

RAF_i = Relative Absorption Factor (inhalation) ET = Duration of each exposure event (hr/day)

EF = Exposure frequency (days/year)

EP = Duration of the exposure period (years)

CF = Appropriate unit conversion factor

BW = Body weight of the receptor of concern during the averaging period

AP = Averaging period

Average Daily Dose for the Respiratory System (ADD_{inhal})

$$ADD_{inhal} = \frac{[C_{part}] \times 0.5 \times [PM_{10}] \times IR_{air} \times RAF_{i} \times ET \times EF \times EP \times CF}{BW_{i} \times AP}$$

where:

ADD_{inhal} = Average Daily Dose due to inhaled particulates entering the lungs (mg/kg-day)

[C_{part}] = Concentration of constituent in airborne particulates (mg/kg)

 $[PM_{10}]$ = Concentration in air of particulates less than 10 microns in diameter ($\mu g/m^3$)

IR_{air} = Inhalation rate (m³/hr) RAF_i = Relative Absorption Factor



ET = Duration of each exposure event (hr/day)

EF = Exposure frequency (days/yr)

EP = Duration of the exposure period (years)

CF = Appropriate unit conversion factor (10^{-9} kg/µg)

BW = Body weight (kg) AP = Averaging period (d)

Dose response values for inhalation exposure (i.e., unit risk factor and reference concentration) are expressed on a mass of compound per volume of air basis, using the assumption that a person weighs 70 kg and has a daily inhalation rate of 20 m³. Therefore, prior to the characterization of risk, ADD_{inhal} (mg/kg-day) for the compound must be converted to an average daily exposure (ADE_{inhal}) (mg/m³) in order to make it compatible with the corresponding dose response values. This can be accomplished using the following equation:

$$ADE_{inhal} = \frac{ADD_{inhal} \times BW_{assumed}}{Inh_{assumed}}$$

where:

 ADE_{inhal} = Average daily exposure COPC concentration (mg/m³)

ADD_{inhal} = Average daily dose due to inhaled particulates in the lungs (mg/kg-d)

BW_{assumed} = Body weight assumed in the development of RfCs and URFs (70 kg)

Inhalation rate assumed in the development of RfCs and URFs (20 m³/d)

For the evaluation of the trespasser, the concentration of PM_{10} in air was assumed to be $32 \,\mu g/m^3$ (residential concentration from MADEP, 2008).

3.3.8.3 Asbestos in Soil

As described in the Phase II CSA, AMEC collected 42 soil samples for asbestos analysis in April 2011. Visual and laboratory identification of asbestos were in agreement, with only one location presenting asbestos. Three aliquots from this location were submitted for elutriator testing to estimate airborne asbestos concentrations. These concentrations are expressed in million fibers per gram of soil (MFG) and were converted to airborne concentrations assuming they comprise the respirable (PM10) fraction of ambient dust:

Where:

OHM_{soil} = Asbestos fiber concentration reported by lab (MFG)

 PM_{10} = Respirable dust concentration in air (mg/m³)



C = Conversion factor
$$\left(\frac{g \text{ soil}}{mg \text{ soil}} \frac{m^3}{ml}\right)$$

Risk is then calculated as follows:

ELCR =
$$ADE_{air} \times IUR$$

Where the average concentration of the oil or hazardous material in air (ADE_{air}) over the exposure period is calculated as

$$ADE_{air} = \frac{OHM_{air} \times EF \times EP}{AP}$$

where:

 OHM_{air} = Concentration of asbestos (f/ml)

EF = Exposure frequency
EP = Exposure period
AP = Averaging period

The risk calculations for asbestos appear in **Attachment D**.

3.3.8.4 Groundwater

Exposure to groundwater used as a residential water supply was assumed to occur via incidental ingestion, dermal contact, and inhalation of vapors while showering. Exposure and risk were estimated using the MassDEP Shortforms. The equations, assumptions and calculations appear in **Attachment E.**

3.3.9 Relative Absorption Factors (RAFs)

The premise of calculating risk or hazard using toxicity data from laboratory experiments is that potential human exposure dose is similar to the administered dose or applied dose in the laboratory experiment. The animal-derived cancer slope factors (CSFs) and reference doses (RfDs) used in quantitative risk assessment were based on applied doses in most cases. However, the efficiency of COPC absorption via a particular route and from a particular matrix (e.g., soil, water) under environmental exposure conditions may differ from the absorption efficiency for the exposure route and matrix used in the experimental study that serves as the basis for the CSF or RfD. RAFs are used to adjust the exposure dose based on these two μ absorption efficiencies. As recommended by MADEP (1995), RAFs for COPCs were derived and used in the calculation of human exposure to soil in the Method 3 evaluation of soil at the DDA.

The RAFs used in the RC were obtained from MADEP (2009) and are shown in **Table 3-5**. A value of 1 is used for inhalation RAFs (including asbestos).



3.4 RISK CHARACTERIZATION

3.4.1 Method 3 Risk Characterization

A Method 3 RC was selected for evaluating potential risk at the DDA. In the Method 3 approach, risk characterization is the step in the risk assessment process that combines the results of the exposure assessment and the toxicity assessment for each COPC in order to estimate the potential for noncarcinogenic and carcinogenic human health effects from exposure to that constituent. This section summarizes the results of the RC for each receptor evaluated in this risk assessment. **Tables 3-7** and **3-8** summarize the total noncarcinogenic risks and carcinogenic risks estimated for each receptor for soil and groundwater, respectively. RC calculations are presented in **Attachment C**.

3.4.1.1 Noncarcinogenic Risk Characterization of Soil

The potential for exposures to COPCs in soil at DDA to result in adverse noncarcinogenic health effects was estimated for each receptor by comparing the Average Daily Dose (ADD) for each constituent (derived in Section 3.3.8) with the Reference Dose for that constituent (presented in Section 3.2.1). The resulting ratio, which is unitless, is known as the Hazard Quotient (HQ) for that constituent. The HQ is calculated using the following formula:

$$HQ = \frac{ADD}{RfD}$$

where:

HQ = Hazard Quotient (unitless);

ADD = Average Daily Dose (mg/kg-day); and

RfD = Reference Dose (mg/kg-day).

When a Hazard Quotient for a COPC does not exceed 1, the Reference Dose has not been exceeded, and no adverse noncarcinogenic health effects are expected to occur as a result of exposure to that COPC via that route. The HQs for each pathway are summed to yield the Hazard Index (HI) for that COPC. A total HI for the receptor is estimated by summing the COPC-specific HIs. A total HI for a receptor that does not exceed 1 indicates that no adverse noncarcinogenic health effects are expected to occur as a result of that receptor's potential exposure to COPCs identified at the Site.

Table 3-7 presents the total HIs calculated for each receptor. The total HIs do not exceed MADEP's noncancer risk limit of 1 for any receptor. Accordingly, DDA soil achieves a



condition of No Significant Risk of harm to human health with respect to noncarcinogenic effects.

3.4.1.2 Carcinogenic Risk Characterization of Soil

The purpose of carcinogenic risk characterization is to estimate the likelihood, over and above the background cancer rate, that a receptor will develop cancer in his or her lifetime as a result of site-related exposures to COPCs in various environmental media. This likelihood is a function of the dose of a constituent and the Cancer Slope Factor (CSF) for that constituent. The relationship between the Excess Lifetime Cancer Risk (ELCR) and the estimated Lifetime Average Daily Dose (LADD) of a constituent may be expressed as:

ELCR = 1- e CSF*LADD

where:

ELCR = Excess Lifetime Cancer Risk (unitless);

CSF = Cancer Slope Factor (1/(mg/kg-day)); and

LADD = Lifetime Average Daily Dose (mg/kg-day).

When the product of the CSF and the LADD is much greater than 1, the ELCR approaches 1 (i.e., 100% probability). When the product is less than 0.01 (1x10⁻²), the equation can be closely approximated by:

ELCR = CSF × LADD

The product of the equations is unitless, and provides an estimate of the potential carcinogenic risk associated with a receptor's exposure to that constituent via that pathway. ELCRs are calculated for each potentially carcinogenic constituent via each exposure pathway. For each receptor, the total ELCR for each COPC is calculated by summing the potential risks derived for each pathway by which the receptor is assumed to be exposed. A Total ELCR for the receptor is then calculated by summing the COPC-specific ELCRs.

ELCRs estimated for all receptors are presented in Table 3-7. Total ELCRs are less than MADEP's cancer risk limit of 1 x 10⁻⁵ for all receptors. Accordingly, DDA soil achieves a condition of No Significant Risk of harm to human health with respect to carcinogenic effects.

3.4.1.3 Risk Characterization of Groundwater as a Potable Supply

Groundwater at the Site is categorized as GW-1 and GW-3. It is unlikely that groundwater at the Site will be used for drinking water. However, because groundwater has been designated as a Potential Drinking Water Source by the Town of Walpole, exposure and risk associated



with use of groundwater as a potable supply (human heath) has been assessed as part of this RC. These estimated risks appear in **Table 3-8** and **Appendix E**..

3.4.2 Summary of Human Health Risk Characterization

Potential noncarcinogenic and carcinogenic risks associated with exposure to COPCs in DDA/SFS soil and groundwater are below 1 and 1 x 10⁻⁵, respectively. **Therefore, soil and groundwater in the DDA and SFS achieve a condition of No Significant Risk**.

3.4.3 Applicable or Suitably Analogous Public Health Standards

The MCP at 310 CMR 40.0993 (3) requires an evaluation of Applicable and Suitably Analogous Standards (ASAS) in addition to quantitative risk characterization. The general list of ASAS provided in the MCP includes the Massachusetts Drinking Water Quality Standards and the Massachusetts Air Quality Standards. These two sets of ASAS are applicable to the Site.

The Massachusetts Air Quality Standards relate to ambient concentrations of the so-called "criteria pollutants" (sulfur oxides, particulate, carbon monoxide, ozone, nitrogen dioxide and lead). None of the COPCs at the Site is a criteria pollutant.

The Massachusetts Maximum Contaminant Levels (MMCLs) at 310 CMR 22.00 were compared with DDA groundwater concentrations, as groundwater is categorized as GW-1. **Table 3-2** presents the comparison to MMCLs. **Concentrations of arsenic in some wells exceed MMCLs.** However, as described in Section 3.1 and Attachment B, arsenic represents a background condition and therefore not pose Significant Risk.

3.4.4 Risk of Harm to Safety

In accordance with 310 CMR 40.0994, the risk of harm to safety was evaluated. No structures are present at the Site. No overt situations posing a threat of physical harm or bodily injury exist. As such, the Site demonstrates a condition of No Significant Risk of Harm to Safety.

3.4.5 Risk of Harm to Public Welfare

In accordance with 310 CMR 40.0994, the risk of harm to public welfare was evaluated. No dangerous or nuisance conditions exist at the Site, nor are persistent odors reported. As part of the public welfare evaluation, EPCs for COPCs in Site soil and groundwater were compared to Upper Concentration Limit (UCLs). **Tables 3-1** through **3-3** shows these comparisons for soil and groundwater. No EPCs exceeded soil UCLs or groundwater UCLs. **Therefore, the Site achieves a condition of No Significant Risk of Harm to Public Welfare.**



3.4.6 Human Health Risk Characterization Conclusions

Conditions at DDA represent a condition of No Significant Risk with respect to human health, safety, and public welfare. Although arsenic concentrations in groundwater exceed MMCLs, arsenic concentrations are consistent with background and therefore do not represent Significant Risk. No other Applicable and Suitably Analogous Standards are exceeded.

3.5 Uncertainty Analysis of the Human Health Risk Characterization

Within any of the four steps of the risk assessment process, assumptions must be made due to a lack of absolute scientific knowledge. Some of the assumptions are supported by considerable scientific evidence, while others have less support. Every assumption introduces some degree of uncertainty into the risk assessment process. Conservative assumptions are made throughout the risk assessment to ensure that the health of local populations and the environment are protected. Therefore, when all of the assumptions are combined, it is much more likely that actual risks, if any, are over-estimated rather than under-estimated.

The assumptions that introduce the greatest amount of uncertainty in this risk assessment are discussed in this section. They are discussed in general terms, because for most of the assumptions there is not enough quantitative information to assign a numerical value that can be factored into the calculation of risk.

3.5.1 Hazard Identification

During the Hazard Identification step, constituents are selected for inclusion in the quantitative risk characterization. COPCs were selected based on potential association with historical Site activities and comparison to background conditions. Detected constituents were screened against background concentrations for both the human health and ecological risk characterizations. Maximum concentrations were also screened against ecological risk-based concentrations for the selection of COPECs for the ecological risk characterization (Section 4.1.2). It is unlikely that constituents have been overlooked in the MCP protocols utilized for the several rounds of sampling and analytical methodologies conducted at the property.

3.5.2 Toxicity Assessment

Dose response values are usually based on limited toxicological data. For this reason, a margin of safety is built into estimates of both carcinogenic and noncarcinogenic risk, and actual risks are lower than those estimated. The two major areas of uncertainty introduced in the dose response assessment are: (1) animal to human extrapolation; and (2) high to low dose extrapolation. These are discussed below.



Human dose response values are often extrapolated, or conservatively estimated, using the results of animal studies. Extrapolation from animals to humans introduces a great deal of uncertainty in the risk assessment because in most instances, it is not known how differently a human may react to the constituent compared to the animal species used to test the constituent. The procedures used to extrapolate from animals to humans involve conservative assumptions and incorporate several uncertainty factors that over-estimate the adverse effects associated with a specific dose. As a result, over-estimation of the potential for adverse effects to humans is more likely than under-estimation.

Predicting potential health effects from the exposure to site soil requires the use of generally-recognized models to extrapolate the observed health effects from the high doses used in laboratory studies to the anticipated human health effects from low doses experienced in the environment. The MCP-specified models contain conservative assumptions to account for the large degree of uncertainty associated with this extrapolation (especially for potential carcinogens) and therefore, tend to be more likely to over-estimate than under-estimate the risks.

3.5.3 Exposure Assessment

During the exposure assessment, average daily doses of COPCs to which receptors are potentially exposed are calculated, which involves assumptions about how often exposure occurs. Such assumptions include location, accessibility, and use of an area. With this in mind, the receptor, or person who may potentially be exposed, and the location of exposure, were both defined for this risk assessment. The locations where certain activities were assumed to take place have been purposely selected because chemical concentrations and frequency of exposure are expected to be high (i.e., use of the maximally affected areas).

Exposures and risks in this RC are based on the assumption of trespassing on a fairly routine basis. In fact, the DDA is not readily accessible and is unlikely to serve as a trespassing destination for local adolescents. In addition, asbestos exposure was based on a single identified location with surficial asbestos, the maximum modeled airborne concentration at that location, and the assumption that all inhalation while trespassing is at that location. These are highly conservative assumptions that are likely to vastly over-estimate the actual degree of fiber inhalation that might be experienced when traversing the area.

3.5.4 Risk Characterization

The risk of adverse human health effects depends on estimated levels of exposure and on dose response relationships. Once exposure to and risk from each of the selected constituents is calculated, the total risk posed by exposure to site soil is determined by combining the health risk contributed by each constituent. Where COPCs do not interact, do not affect the same target organ or do not have the same mechanism of action, summing the risks for multiple



COPCs results in an over-estimate of risk posed by the Site. However, in order not to understate the risk, it is assumed that the effects of different constituents may be added together. Overall, this conservative method of risk characterization is expected to over-estimate, rather than under-estimate, health risks posed by the Site. Because all potentially complete exposure pathways were evaluated in the HHRC, potential risks are not likely to be underestimated.



4.0 RISK TO THE ENVIRONMENT

Risk to the Environment was evaluated in this RC in accordance with MADEP guidance *Method 3 Environmental Risk Characterization* (MADEP, 1996). This MADEP guidance provides for two stages of environmental risk characterization:

- 1. Stage I Environmental Screening, which is used to identify those situations which require further evaluation; and
- 2. Stage II Environmental Risk Characterization, which is a detailed evaluation of those environmental exposure pathways identified in Stage I.

As part of Stage I, available site information is reviewed to identify the presence of environmental receptors and to determine whether the identified receptors are currently exposed, or could potentially be exposed, to site-related constituents. According to MADEP guidance, a "complete exposure pathway means that the contamination is actually reaching plants or animals, or is likely to do so in the future" (MADEP, 1996). Exposure pathways that are not complete and are not likely to be complete, are not required to be evaluated further.

As discussed in Section 3.1.1, the DDA is limited to upland areas and does not include Cedar Swamp Brook. Therefore, this ERC is limited to evaluation of surface soil.

4.1 Stage I Environmental Screening

Information about the Site was reviewed to determine whether environmental receptors are currently or may in the future be present, and to determine whether the identified receptors are currently exposed, or could potentially be exposed, to Site-related constituents of potential ecological concern (COPECs). Environmental screening begins with an assessment of whether the potential for exposure exists at the Site. Exposure pathways, or links between the presence of a constituent and ecological receptors, must be complete for exposure potential to exist.

4.1.1 Habitat Evaluation

AMEC conducted field evaluations of natural habitats in several areas of the BMC property to provide specific information about ecological communities and wildlife present in the area. These observations were used to support the ecological risk characterization and select representative species to be used in Stage II environmental RC. A field inspection was conducted on September 15, 2009, that documented species present in the DDA area. A field visit performed in August 2011 as part of the asbestos soil sampling provides additional, more recent, documentation as to current conditions.

Much of the former BMC property is disturbed by historical industrial activity. The property contains roads and other corridors (cleared paths, railroad spur), impervious areas, buildings, former building locations, areas of stockpiled soil, rock, construction debris, and channelized



water bodies. The property is surrounded by both undeveloped land (forested uplands and wetlands) and developed areas. Land uses in developed areas include residential, commercial, and industrial. The Neponset River flows from southwest to northeast through the eastern portion of the property; it has been channelized where it crosses the Site, and a part of the river is diverted into Ruckaduck Pond. The tributary Cedar Swamp Brook approaches the northern border of the Site from the west, and joins the Neponset east of the Site in a forested wetland system. A railroad right-of-way forms part of the western border of the Site. These features are important factors influencing the vegetative communities and the wildlife species that use the Site. Most areas of the Site are vegetated at least in part by invasive species and early colonizing species of disturbed lands.

Prior to the field evaluation, AMEC reviewed existing project documentation, publicly available maps, and related ecosystem information. The habitat evaluation then consisted of observing and describing ecological communities at the Site, classifying these ecological communities in relation to an existing classification system, investigating the Site for evidence of wildlife use of the areas, and based on literature references, identifying wildlife species that could reasonably be expected to occupy the areas of interest.

The reference ecological community description used for the habitat evaluation is the Classification of Natural Communities developed by the NHESP (Swain and Kearsley 2001). This classification system describes ecological communities according to vegetation species composition, landscape, and hydrology in particular environments. The classes are based on exemplary conditions of recurring assemblages of plants and animals as observed in undisturbed landscapes, and generally exclude those areas created or maintained by human activities. Although the NHESP classes may not strictly correspond to the areas under investigation at the Site as a result of historical industrial activities, it provides a convenient reference point for ecosystem description and development of species lists. Detailed field observations of environmental conditions were used with Site personnel reports and literature reports of similar ecosystems, species assemblages, and wildlife habitats to complete the ecological habitat characterization of the areas.

An AMEC ecologist inspected the Site on September 15, 2009, to view the entirety of the Site and the surroundings and describe the ecological communities. Communities were described by the plant strata present (tree canopy, subcanopy, shrubs, herbaceous vegetation, and vines), the vegetation species composition in each, and by a description of landscape including soils and hydrology. The inspection also looked for evidence of recent disturbance or stressed vegetation. Information on wildlife use of the areas was gathered through direct observation of species or their signs, and by comparing the observed habitats with published lists of animals known or expected to occur in those habitats.

The DDA is a highly disturbed area. The majority of the forested upland communities within the project area are unexceptional white pine / oak forests. These communities are dominated by



red oak and white pine in the canopy. The understory of this community is in large part dominated by oriental bittersweet, glossy buckthorn, and poison ivy. Debris from old machines and construction activities are present in this community.

Cover is characterized by a mix of grasses, shrubs and gravelly zones, with a road running through and areas of standing water. Habitat is of overall limited quality due to large patches of unvegetated surface. Trees are limited to the perimeter. Many of the flora species observed in the DDA area (listed below) are invasive species representative of highly disturbed areas:

Canopy - Oak Forest

Quercus rubra Red oak Acer rubrum Red maple White pine Pinus strobusstrobes White oak Quercus alba Quaking aspen Populus tremula Fraxinus americana White ash Ulmus americana American elm Betula alleghaniensis Yellow birch Hemlock species

Shrub Layer

Elaeagnus angustifolia Russian olive Frangula alnus Glossy buckthorn

Myric spp. Bayberry Betula populifolia Gray birch Toxicodendron radicans Poison ivv

Oriental bittersweet Celastrus orbiculatus Rosa multiflora

Multiflora rose

Grape

Understory

Toxicodendron radicans Poison ivy Parthenocissus quinquefolia Virginia creeper

Arctium spp Burdock Graminoides Grasses

Meadow

Ambrosia artemisifolia Common ragweed Goldenrod spp. Solidago and Euthamia spp. Daucus carota Queen Anne's lace Populus spp Tree seedlings Fallopia japonica Japanese knotweed

Trifolium spp. Red clover

Rudbeckia hirta Black-eyed susan

Scrophulariaceae spp. Mulleins



Wildlife observed or expected consists of those species normally inhabiting disturbed areas within and near forest edges, and mid-late old field successional communities, including suburban areas, roadsides, and waste areas. Deer tracks, and burrows were noted, as well as frogs and crickets. The following species have been observed or would be expected in the DDA area:

Amphibians

Bufo americanus American toad Bufo woodhouseii fowleri Fowler's toad

Rana pipiens Northern leopard frog

Reptiles

Chelydra serpentina Snapping turtle Chrysemas picta Snapping turtle

Coluber constrictor Northern black racer
Diadophis punctatus edwardsii Northern ringneck snake
Lampropeltis triangulum Eastern milk snake

Lichlorophis vernalis Eastern smooth green snake

Storeria dekayi Northern brown snake
Storeria occipitomaculata Northern redbelly snake
Thamnophis sirtalis Eastern garter snake

Avian Species

Baeolophus bicolorTufted titmouseBombucilla cedrorumCedar waxwingBubo virginianusGreat-horned owlButeo jamaicensisRed-tailed hawkCardinalis cardinalisNorthern cardinalCarduelis tristisAmerican goldfinch

Carpodacus mexicanus House finch
Cathartes aura Turkey vulture
Corvus brachvrhvnchos American crow

Cyanocitta cristata

Melospiza melodia

Passer domesticus

Turdus migratorius

Zenaida macroura

American crow

Blue jay

Song sparrow

House sparrow

American robin

Mourning dove

<u>Mammals</u>

Blarina brevicauda Northern short-tailed shrew

Canis latrans Coyote
Didelphis marsupialis Opossum

Didelphis virginianaVirginia opossumMicrotis pennsylvanicusMeadow voleOdocoileus virginianusWhite-tailed deer



Mammals

Peromyscus leucopus White-footed mouse

Procyon lotorRaccoonScalopus aquaticusEastern moleSciurus carolinensisGray squirrel

Sylvilaqus floridanus Eastern cottontail rabbit Tamias striatus Eastern chipmunk

4.1.2 Comparison of Soil Concentrations to Ecological Risk-Based Concentrations

As required in 310 CMR 40.0995, maximum detected concentrations in soil from 0 to 2 ft bgs in the DDA were compared to MADEP's background concentrations from "natural" soil (MADEP, 2002a) and ecological risk-based concentrations in soil.

Table 4-1 presents the ecological screening for DDA soil. DDA maximum detected concentrations exceed screening levels for arsenic, several metals, several PAH compounds, and dioxin/furan compounds (TCDD TEQs). No screening criteria exist for the petroleum hydrocarbon fraction, so these have been included as COPECs for DDA. Because the Stage I screening process identified COPECs, a Stage II ERC was performed.

4.2 Stage II Environmental Risk Characterization

Environmental risk characterizations involve multiple steps. The first step, Problem Formulation, develops a Conceptual Site Model (CSM) that identifies ecological resources (Site biota) and identified the assessment endpoints (ecological "entities" and their characteristics and functions targeted for protection). Based on the key assessment endpoints, measurement endpoints (quantitative or measurable characteristics or attributes of the assessment endpoints) are developed, such as measures or estimates of exposure or effect. Risk characterization, the third step, evaluates the likelihood of adverse ecological effects on the assessment endpoints based on the results of the analysis. Each of these steps is described in more detail below.

4.2.1 Problem Formulation

Problem formulation includes the following:

- 1. Review of available data on ecological communities and existing data on constituent concentrations in environmental media (accomplished during the Stage I Environmental Screening);
- 2. Development of a conceptual site model (CSM);
- 3. Identifying assessment endpoints;
- 4. Selection of representative ecological receptor species; and
- 5. Selection of COPECs.



4.2.1.1 Development of a Conceptual Site Model

One important component of the Problem Formulation phase is the development of a CSM, which integrates existing knowledge of the physical, biological, and constituent conditions at the Site into a strategy for assessing whether concentrations of COPECs in soil or prey pose a potential threat to exposed populations.

CSMs describe the direct and indirect pathways through which ecological receptors might be exposed to COPECs. According to U.S. EPA (1998), the CSM presents a verbal description and visual representation of the pathways from constituent sources to potentially exposed receptors. The two objectives of a CSM are to: (1) illustrate the important relationships within the ecosystem and (2) specify exposure scenarios to be evaluated in the environmental risk characterization. The CSM for the Site is presented in **Figure 4-1**.

The CSM suggests ecological receptors may be directly exposed to COPECs through contact with soil and indirectly exposed by consumption of food organisms that have accumulated COPECs in their tissues due to soil exposure.

4.2.1.2 Assessment and Measurement Endpoints

Assessment endpoints represent an explicit expression of the actual environmental values to be protected at the Site. Assessment endpoints are either measured directly or evaluated through indirect measures. They have been developed based on the four selection criteria in the U.S. EPA guidance (1997a) and the requirement that the ecosystems, communities, and/or species selected as endpoints are present at the Site. The four selection criteria include:

- 1. Consideration of the constituents present and their concentrations;
- 2. Mechanisms of toxicity of the constituents to different groups of organisms;
- 3. Ecological relevance of receptor groups that are potentially sensitive or highly exposed to the constituent and attributes of their natural history; and
- 4. Potential completeness of exposure pathways from the constituent(s) to the relevant receptor group.

Measurement endpoints represent quantifiable ecological characteristics that can be measured, interpreted, and related to the valued ecological component(s) chosen as the assessment endpoints. When selecting measurement endpoints, there should be an explicit relationship between the measurement endpoint and the assessment endpoint (i.e., value to be protected) to which it is linked. Measurement endpoints should be selected based on the species/community/habitat patterns across the Site, the varying relationship to the COPEC concentrations, and considerations of the mechanisms of toxicity.

Assessment and measurement endpoints have been developed to evaluate COPEC concentrations and potential ecological risks at the Site. In some cases, more than one



measurement endpoint may be identified for a particular assessment endpoint. These instances permit a "weight-of-evidence" approach to be used in risk characterization, in which a measurement endpoint that has more relevance to the assessment endpoint than another measurement endpoint is assigned more "weight" in the interpretation of risk estimates. In other cases, a measurement endpoint may be relevant to more than one assessment endpoint, or a secondary assessment endpoint may not have a measurement endpoint associated with it

Based on the CSM for the Site (**Figure 4-1**) consisting of primary producers, primary consumers (herbivores), and secondary consumers (insectivores and carnivores), the following assessment and measurement endpoints have been evaluated for each of the three exposure areas:

<u>Assessment Endpoint 1</u> – The reproductive success and population sustainability of herbivorous mammal and bird populations potentially exposed to COPECs in surface soil and prey.

<u>Corresponding Measurement Endpoint</u> – Comparison of predicted average daily doses of COPECs to toxicity reference values (TRVs) for mammalian and avian receptors.

<u>Assessment Endpoint 2</u> – The reproductive success and population sustainability of insectivorous mammal and bird populations potentially exposed to COPECs in surface soil and prey.

<u>Corresponding Measurement Endpoint</u> – Comparison of predicted average daily doses of COPECs to TRVs for mammalian and avian receptors.

<u>Assessment Endpoint 3</u> – The reproductive success and population sustainability of carnivorous mammal and bird populations potentially exposed to COPECs in surface soil and/or prey.

<u>Corresponding Measurement Endpoint</u> – Comparison of predicted average daily doses of COPECs to TRVs for mammalian and avian receptors.

4.2.1.3 Selection of Ecological Receptors

Criteria for the selection of wildlife receptors include three factors specified in U.S. EPA guidance (1989, 1992, 1997c) for determining "key organisms" in an ecological food web: (1) resident communities or species exposed to constituent concentrations in soil and surface water; (2) species or functional groups considered to be essential to, or indicative of, the normal functioning of the affected habitat; and (3) species representing federal or state threatened or endangered species. Other factors to be considered include the trophic level of the organisms, area use factors and feeding habits of the species, availability of life history and toxicity data for the species, and abundance of the species.



According the U.S. Fish and Wildlife Service published data "Federally Listed Endangered and Threatened Species in Massachusetts" revised 6/22/2009, (http://www.fws.gov/newengland/) there are no known occurrences of federally-listed species in Walpole, Massachusetts.

AMEC submitted a written request to the Massachusetts Division of Fisheries and Wildlife (DFW) Natural Heritage and Endangered Species Program (NHESP) for information regarding state-listed endangered, threatened, or species of special concern that may be known by NHESP to inhabit the vicinity of the site. NHESP responded with a letter on October 28, 2009 (NHESP, 2009) stating that, based on the information provided, the Site, or a portion thereof, is located within Priority Habitat 1072 (PH 1072) and Estimated Habitat 855 (EH 855) as indicated in the Massachusetts Natural Heritage Atlas (13th Edition). NHESP's database indicates that one state-listed rare species has been found in the vicinity of the Site: Callophrys hesseli (Hessel's Hairstreak), a Species of Special Concern. Hessel's hairstreak is a butterfly (Lepidoptera) that is protected under the Massachusetts Endangered Species Act (MESA) (M.G.L. c. 131A) and its implementing regulations (321 CMR 10.00). State-listed wildlife are also protected under the state's Wetlands Protection Act (WPA) (M.G.L. c. 131, s. 40) and its implementing regulations (310 CMR 10.00). Hessel's Hairstreak exclusively inhabits Atlantic white cedar (Chamaecyparis thyoides) swamps and bogs. No Atlantic white cedar swamps or bogs exist at DDA. As such, it is highly unlikely that the species Hessel's hairstreak occurs in the investigated areas.

The following representative receptors have been selected for evaluation of the measurement endpoints discussed above.

<u>Herbivorous Mammals.</u> The meadow vole (*Microtus pennsylvanicus*) is a small plant-eating mammal commonly found in vegetated fields. They may also incidentally ingest surficial soils while feeding.

<u>Herbivorous Birds.</u> The quail, also referred to as northern bobwhite, (*Colinus virginianus*) subsists mainly on seed and low-lying vegetation. They are commonly found in grassy fields and pastures. They may also incidentally ingest surficial soils while feeding.

<u>Insectivorous Mammals.</u> The short-tailed shrew (*Blarina brevicauda*) is a species that preys mainly on insects and soil invertebrates. They can be found in almost any habitat, but prefer habitats with vegetation. They may also incidentally ingest surficial soils while feeding.

<u>Insectivorous Birds.</u> The American woodcock (*Scolopax minor*) subsist primarily on invertebrates. They are commonly found in woodlands and open fields. They may also incidentally ingest surficial soils while feeding.



<u>Carnivorous Mammals.</u> The red fox (*Vulpes vulpes*) primarily feeds on small animals. They can be found in a large variety of habitats, including abandoned fields and woodlands. They may also incidentally ingest surficial soils while feeding.

<u>Carnivorous Birds.</u> The red-tailed hawk (*Buteo jamaicensis*) subsists primarily on small mammals and can be found in woodlands and open fields. They may also incidentally ingest surficial soils while feeding.

4.2.1.4 Selection of COPECs

Constituents with maximum detected soil concentrations exceeding the screening concentrations presented in Section 4.1.2 were selected as COPECs. COPECs are shown in **Table 4-1** for DDA. Soil is the only identified medium of ecological concern

4.2.2 Analysis

The analysis stage of environmental RC consists of two steps: the exposure assessment and the effects assessment. These are described in the sections below.

4.2.2.1 Exposure Assessment

The exposure assessment estimates the magnitude, frequency, duration, and types of potential exposures to COPECs in food webs at the Site. This includes calculating EPCs in Site media and identifying equations and exposure parameter values used to estimate potential exposure for the ecological receptors. Potential receptors and exposure routes evaluated in the exposure assessment were identified in Section 4.2.1 above. The following sections describe the equations and exposure assumptions used to estimate potential exposure and derivation of EPCs. Exposure parameter values are presented in **Table 4-2**.

A food chain analysis was used to estimate exposure to Site COPECs for each representative receptor. The general calculation is

$$ADD\left(\frac{mg}{kg-day}\right) = \frac{\left[\left(C_f x I R_f\right) + \left(C_s x I R_s\right)\right] x AUF}{BW}$$

where:

ADD = Average daily dose (mg/kg-day)

C_f = Concentration of COPEC in food (mg/kg)

 IR_f = Ingestion rate of food (kg/day)

 C_s = Concentration of COPEC in soil (mg/kg) IR_s = Incidental ingestion rate of soil (kg/day)



AUF = Area use factor (unitless) and

BW = Body weight (kg).

The parameter values used in the above equation for each receptor, along with the reference, are shown in **Table 4-2**. Detailed calculations appear in **Attachment F**.

4.2.2.2 Exposure Point Concentrations

Using the methodology described in Section 3.3.7.1, soil EPCs were calculated as the arithmetic mean of concentrations from samples in 0 to 3 foot depth interval. Soil EPCs for ecological risk calculations are presented in **Table 4-1** for DDA.

Earthworm samples were collected from the DDA area. One sample location falls within the DDA boundary and was used to represent invertebrate dietary concentrations. For COPECs with no earthworm analytical results, the COPEC concentration in invertebrates was estimated by multiplying EPCs in soil by a constituent-specific soil-to-soil invertebrate BTF obtained from U.S. EPA (1999, 2007; see **Table 4-3**). For some inorganics, tissue concentrations are calculated using published logarithmic equations that reflect declining BTFs with increasing soil concentrations. These calculations appear in **Attachment F.**

4.2.3 Effects Assessment

The effects assessment entails a review of the ecotoxicology of the COPECs and development of TRVs for the selected ecological receptors for each COPEC. TRVs were selected using toxicity information developed by the U.S. EPA for the derivation of Ecological Soil Screening Levels (EcoSSLs) and toxicity information from other sources such as Oak Ridge National Laboratory (Sample et al., 1996).

Chronic No Observable Adverse Effect Level (NOAEL) and Lowest Observable Adverse Effect Level (LOAEL) values for mammals and birds were taken from literature sources such as Schafer et al. (1983); and databases including those available from the Oak Ridge National Laboratory (Sample et al., 1996), and the U.S. EPA Integrated Risk Information System (IRIS; U.S. EPA, 2010a). The selection of appropriate TRVs focused on identifying NOAELs and LOAELs for the following toxic effect endpoints in the following order of preference: (1) reproduction, (2) growth/development, and (3) survival. The lowest TRV for a particular constituent was selected from the sources cited above. Both NOAEL-based TRVs and LOAEL-based TRVs were used in this risk characterization.

For constituents without chronic NOAELs, but for which other toxicity values were available, uncertainty factors were applied to estimate chronic NOAELs from the available toxicity data. These other toxicity values include less-than-chronic NOAELs (e.g., subchronic NOAELs), LOAELs, and the lethal dose for 50 percent of a study population (LD₅₀). An uncertainty factor of 10 (as cited in Sample et al., 1996) was used to adjust subchronic LOAEL TRVs to NOAEL



TRVs, and an uncertainty factor of 10 (as cited in Sample et al., 1996) was used to adjust subchronic TRVs to chronic TRVs. The Standard Practice for Wildlife Toxicity Reference Values (USACHPPM, 2000) recommends the use of a total uncertainty factor of 100 to adjust LD_{50} values to chronic NOAEL equivalent values.

If no toxicity values were available for a particular constituent, the chronic NOAEL or LOAEL for an appropriate surrogate constituent was used as the mammalian or avian TRV. Surrogate constituents were selected based on structural chemistry, specifically, the active moiety/functional group of the constituent. TRVs are presented in **Table 4-4** for mammalian receptors and **Table 4-5** for avian receptors.

Toxicity studies for a COPEC may have been conducted on species other than the receptor species evaluated in the ERC. If toxicity values were not based on data for the receptor species, an allometric conversion based on body size (i.e., weight and surface area) was used to extrapolate between species. For the mammalian receptors, the body size-adjusted TRVs, referred to here as "adjusted NOAEL (or LOAEL)-equivalent TRVs," were calculated using the allometric conversion in the following equation (Sample et al., 1996, 1997):

$$TRV_{adj} = TRV_t (BW_t/BW_r)^{(1-SF)}$$

where:

 TRV_{adj} = Adjusted NOAEL (or LOAEL)-equivalent TRV (mg/kg of body weight per day)

 TRV_t = NOAEL (or LOAEL)-equivalent toxicity reference value for test organism

(mg of constituent/kg of body weight per day)

BW_t = Body weight for test organism (kg)
BW_r = Body weight for receptor species (kg)
SF = Body size scaling factor (unitless)

A body size scaling factor (SF) of 0.75 was used in the above equation to extrapolate TRVs between mammalian species (Sample et al. 1996, 1997). A SF of 1 was used in this equation to extrapolate TRVs between avian species. Mineau et al. (1996) identified a mean SF of 1.15 for birds. However, Sample et al. (1996) report that SFs for a majority of the constituents evaluated (29 of 37) by Mineau et al. (1996) were not significantly different from 1. Therefore, a SF of 1 for TRV extrapolation between avian species was determined to be more appropriate.

4.3 Risk Characterization

Risk characterization is the culmination of the preceding steps of the ecological risk assessment and involves three principal components: (1) risk estimation, (2) risk description, and (3) uncertainty analysis. In this step, the risks associated with estimated exposures are characterized, and the strengths, weaknesses, and assumptions employed in the risk



assessment are fully described. In the risk estimation, the exposure assessment and effects assessment profiles from the Analysis phase are integrated to predict the likelihood of adverse effects to receptors.

4.3.1 Risk Estimation

The risk estimation component for the food chain analysis provides a quantitative evaluation of the exposure assessment and effects assessment results. Potential risks to the ecological receptors were estimated using the hazard quotient (HQ) method. In this method, the estimated exposure (the ADD) is compared to the TRV using the following equation:

HQ = ADD/TRV

where:

HQ = Hazard quotient (unitless)

ADD = Average daily dose (output of food web model) (mg/kg-day), and TRV = Toxicity reference value (NOAEL and LOAEL-based) (mg/kg-day).

HQs are calculated separately for each COEPC in each exposure area for each assessment endpoint below. When the HQ does not exceed 1, the estimated potential exposure does not exceed the TRV, indicating that adverse effects are not likely to occur. When the HQ is greater than 1, the estimated potential exposure exceeds the TRV, and adverse effects cannot automatically be ruled out. However, the food chain model contains multiple conservative assumptions that produce a high bias of risk. Therefore, HQs over 1 need to be evaluated to determine if there is really a likelihood of hazard.

HQs for the DDA appear in **Table 4-6**. NOAEL-based HQs did not exceed 1 with the exception of the insectivorous receptors (shrew and woodcock). The maximum HQ was 9 for TCDD TEQs (shrew); other NOAEL-based HQs over 1 were for chromium (2 for shrew and 3 for woodcock), TCDD TEQs for woodcock (2) and vanadium for woodcock (2) The LOAEL-based HQs were all below 1.

These isolated results for receptors associated with one type of diet do not in and of themselves indicate overall hazard to the environment. Calculated insectivorous HQs are frequently over 1 because of the conservative assumptions associated with estimating dietary contaminant concentrations. For TEQs, the invertebrate concentration was estimated using a BTF of 1.5, which was derived from a single study (Meyn et al, 1997). This study was based on applied sludge, which could overestimate bioavailability from more weathered sources. For chromium, the BTF was based on earthworm intake using a single earthworm sample, which was the highest of the three worm samples analyzed. In addition, earthworms represent a relatively



small proportion of the shrew diet, which is 97% insects (Meyn et al, 1997). Overall, therefore, the dietary exposure to these insectivorous species is characterized by considerable uncertainty.

4.3.2 Risk Description

The risk description component of the ERC phase includes: (1) a summary of the risk estimate(s); (2) a discussion of the evidence supporting the risk estimate(s) - i.e., weight of evidence evaluation; and (3) an interpretation of the ecological significance and relevance of the estimate(s).

4.3.2.1 Summary of Risk Estimates

4.3.2.2 Weight of Evidence

The weight-of-evidence is a crucial element of the interpretation of the ERC results, and it is integral to the risk management evaluation. The following factors are some of the key considerations in the weight-of-evidence evaluation of the various risk estimates:

- 1. The relevance of the evidence to the assessment endpoint. The avian and mammalian indicator species chosen as measurement endpoints (meadow vole, quail, short tailed shrew, American woodcock, red fox, and red-tailed hawk) were selected for several reasons. First, all six species selected represent upper (consumer rather than producer) trophic level predators so that potential food chain effects would be considered. Second, these receptors represent major animal families (mammals and birds) and are indigenous to the habitat at the Site. Thus, the risk estimates for the selected receptors are relevant to the ecological values articulated in the assessment endpoints.
- 2. The relevance of the evidence to the CSM describing the physical fate and transport processes and their direct relevance to the assessment endpoints. The CSM and understanding of historical chemical fate and transport at the Site suggest that the concentrations of COPECs in soil are likely due to historical operations. Media samples, receptors, and exposure pathways were selected for evaluation in the ERC based on the information in the CSM. Thus, the resulting risk estimates are relevant to the CSM.
- 3. The confidence in the risk estimate or other information. The confidence and representativeness in the risk estimates is reflected in the selection of the receptors exposure parameter values, BCFs and TRVs. The ERC evaluated receptors that are typically present in habitats like the exposure areas at the Site. The exposure and toxicity values used in the ERC were obtained from MADEP and U.S. EPA sources. As such, there is high confidence that the estimated HQs do not under-estimate potential risk.



4.3.2.3 Ecological Significance and Relevance

The determination of the ecological significance of potential adverse effects to wildlife species from exposures to COPECs included consideration of the following four items:

- 1. The intensity, or severity, of the predicted adverse effect. The HQ approach was used to assess potential risk. When the HQ is greater than 1, the estimated potential exposure exceeds the TRV and a potential risk may exist for individual organisms. However, an HQ exceeding 1 does not indicate that adverse effects are occurring or will occur, even for individual organisms because of the conservatism in the ERC inputs and methods. Moreover, an HQ greater than 1 (particularly, if less than 10) is unlikely to indicate an adverse effect to the population evaluated in the risk characterization.
- 2. The size of the affected area that can be attributed to COPEC-induced unacceptable effects. The surface soil samples used in the RC were collected from locations where historical disposal activities have occurred. There are large areas on the Site where industrial and disposal activities did not occur and where concentrations are likely lower than those used to estimate EPCs and risks. Because more samples were collected from areas expected to have higher concentrations, and fewer samples from large areas where concentrations are not expected to be elevated, any HQs exceeding 1 likely indicate that any area with adverse effects is limited in size.
- 3. Temporal variation and frequency in the occurrence of unacceptable effects. There are several conservative assumptions associated with the estimation of hazard. For metals, the dietary exposure was based on one worm sample, which may not be representative of the whole DDA exposure area nor of the actual diet, which is primarily insects for the shrew. Therefore, it is likely that the ERC over-estimated the exposure for the shrew.
- 4. The capability of the affected area to recover naturally to partial or full recolonization of populations or communities and conditions that existed prior to the introduction of COPECs. There is no indication that concentrations of COPECs in soil have resulted in actual adverse effects to either individual or populations of ecological receptors. Accordingly, "recovery" does not appear necessary. Given the presence of the remnants of the former manufacturing buildings and structures, a large portion of the MBA does not provide suitable habitat for ecological receptors. As such, receptors are not likely to routinely or frequently forage in these areas. Combined with the likely overestimate of actual average concentrations as a result of a biased sampling plan, it is unlikely that adverse effects would occur frequently, if at all.



4.4 Uncertainty Analysis of the Environmental Risk Characterization

There are a number of sources of uncertainty in ecological risk assessments, which can be broadly grouped into three categories: conceptual model uncertainty, parameter values, and model error. Each of these is discussed below.

<u>Conceptual Model Uncertainty</u>: The CSM summarized the fate and transport processes that are believed to have resulted in current conditions at the Site, and have formed the basis for the field investigations, the exposure pathways that were assessed, the receptors of concern, and the assessment and measurement endpoints. There is uncertainty in the completeness of the fate and transport processes and the extent to which they contribute to the potential exposure pathways and the receptors evaluated.

<u>Parameter Values</u>: Because of inherent biological variability and differences in study design, there is some uncertainty associated with the exposure assumptions that were used for dose calculations, biotransfer factors used to estimate EPCs in dietary components, and the TRVs that were used to estimate the risks. Because the values chosen for these variables are conservative, the resulting exposures and risks are unlikely to be underestimated.

Exposure Point Concentrations: Because soil sampling focused on areas where concentrations were expected to be elevated, the EPCs estimated from this dataset likely over-estimate actual average concentrations. Moreover, samples were collected from areas outside the boundaries of the original RTNs were excluded from the dataset. Therefore the EPCs are likely an over-estimate of the soil contacted by terrestrial ecological receptors.

<u>Conservative exposure assumptions</u>: Many of the exposure assumptions were based on field studies performed in other areas but were assumed to be representative of the behavior of these receptors at the Site. The most important of the exposure assumptions was the conservative assumption of a bioavailability factor of 1. The simplified food-chain exposure models used conservative assumptions and it was assumed that the COPECs were 100 percent bioavailable from the exposure media. Because bioavailability virtually always less than 100 percent, the exposures estimated in this ERC likely over-estimate actual exposures.

<u>TRVs</u>: Chronic TRVs were derived from the lowest reported NOAEL or LOAEL of test organisms after applying uncertainty factors to estimate chronic TRVs for the receptors of interest. The uncertainty factors used in TRV development are well established (e.g., U.S. EPA, 1997d). The TRVs identified in this ERC are for screening purposes and do not indicate concentrations levels associated with specific hazards.

<u>Population Risk Estimates</u>. The hazard quotient approach used in this evaluation is based on a sensitive individual receptor. Because conservative exposure assumptions were combined with conservative toxicity assumptions, the resulting risk estimates represent potential effects to



highly exposed, sensitive individuals within the population. As such, the estimated risks all but certainly overestimate potential effects to the populations.

<u>Model Error</u>: There is model error uncertainty in the method used to derive indirect (food-web) uptake. Although the models are based on established fate and transport processes, dietary preferences, and receptor characteristics, they are generic and may not be representative of the processes that may be occurring at the Site.



5.0 SUMMARY AND CONCLUSIONS

In accordance with the requirements of 310 CMR 40.0000 Subpart I of the MCP, a Method 3 RC of harm to human health, public welfare, safety, and the environment has been completed. The Method 3 human health RC of soil and groundwater in the DDA evaluated potential exposures of current and future trespassers, and future hypothetical groundwater use as a potable supply.

Data from site assessment activities as well as information from other sources (e.g., MADEP and U.S. EPA guidance documents and databases) were used to conduct the risk characterization.

The results of the human health risk characterization indicate that a condition of No Significant Risk can be demonstrated for soil and groundwater at the DDA.

The results of the evaluation of risk of harm to safety and public welfare indicates that no unsafe or nuisance conditions exist at the Site. Soil and groundwater constituent concentrations are less than their respective UCLs. As such, a condition of No Significant Risk to public welfare and a condition of No Significant Risk to safety can be demonstrated at the DDA.

The evaluation of potential risk of harm to the environment included a Stage I screening evaluation of the presence of ecological receptors and potential habitat for terrestrial ecological receptors. Potential exposures of herbivorous mammals, herbivorous avians, insectivorous mammals, insectivorous avians, carnivorous mammals, and carnivorous avians to COPECs in soil and the food web were evaluated using a hazard quotient approach.

Hazard quotients were below 1 for all receptors except the short-tailed shrew. For the shrew, the NOAEL HQs exceeding 1 in DDA were below 10 and only for two COCs. These exposures are not expected to cause adverse environmental impacts to short-tailed shrew populations or populations at the Site the because the hazard quotient approach used in this evaluation is based on a sensitive individual receptor. Conservative exposure assumptions are combined with conservative toxicity assumptions, so that the resulting risk estimates overestimate potential effects to the populations. HQs within an order of magnitude of 1 are not likely to be associated with population effects.

Based on the finding that population-level effects are not expected for all receptors, a condition of No Significant Risk of harm to the environment exists at the DDA.

This human health and environmental risk characterization concludes that DDA achieves a condition of No Significant Risk.



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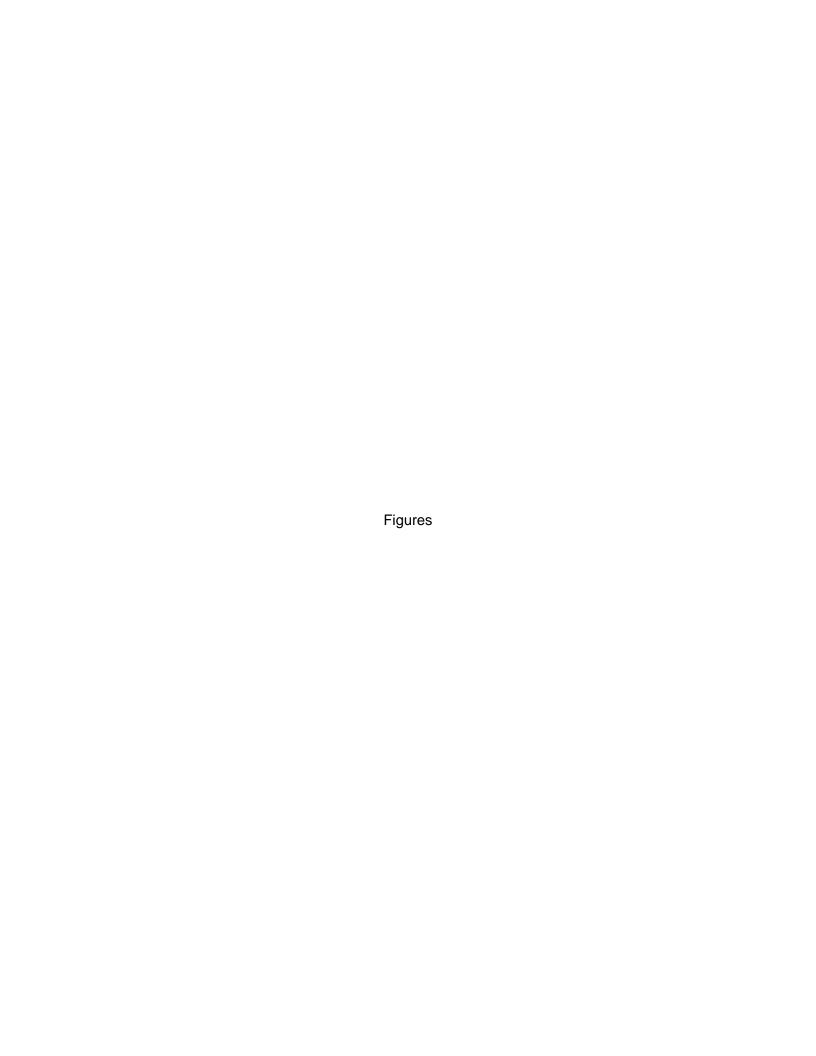
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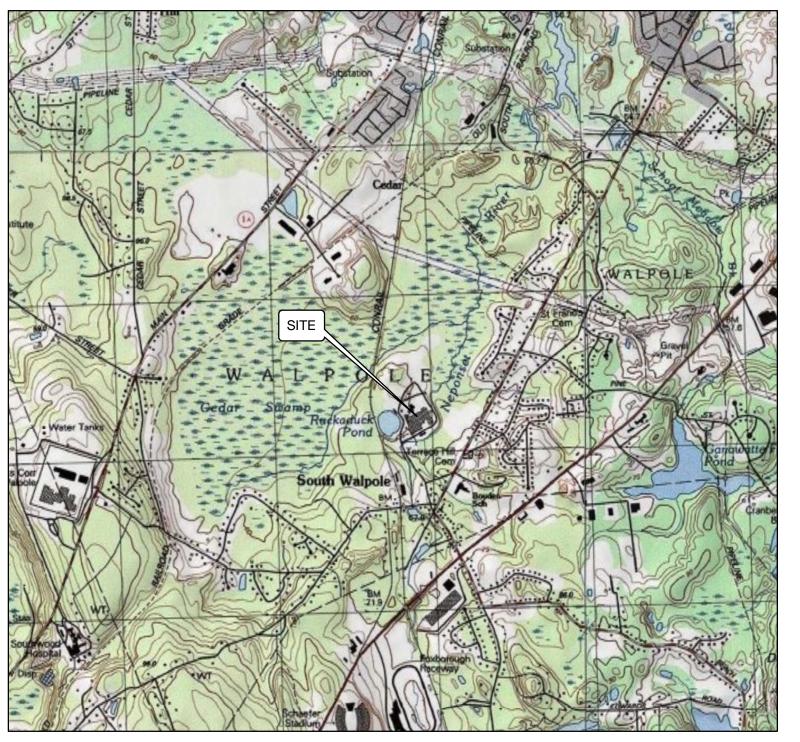


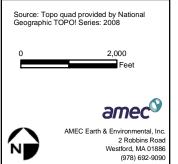
FIGURE 1-1

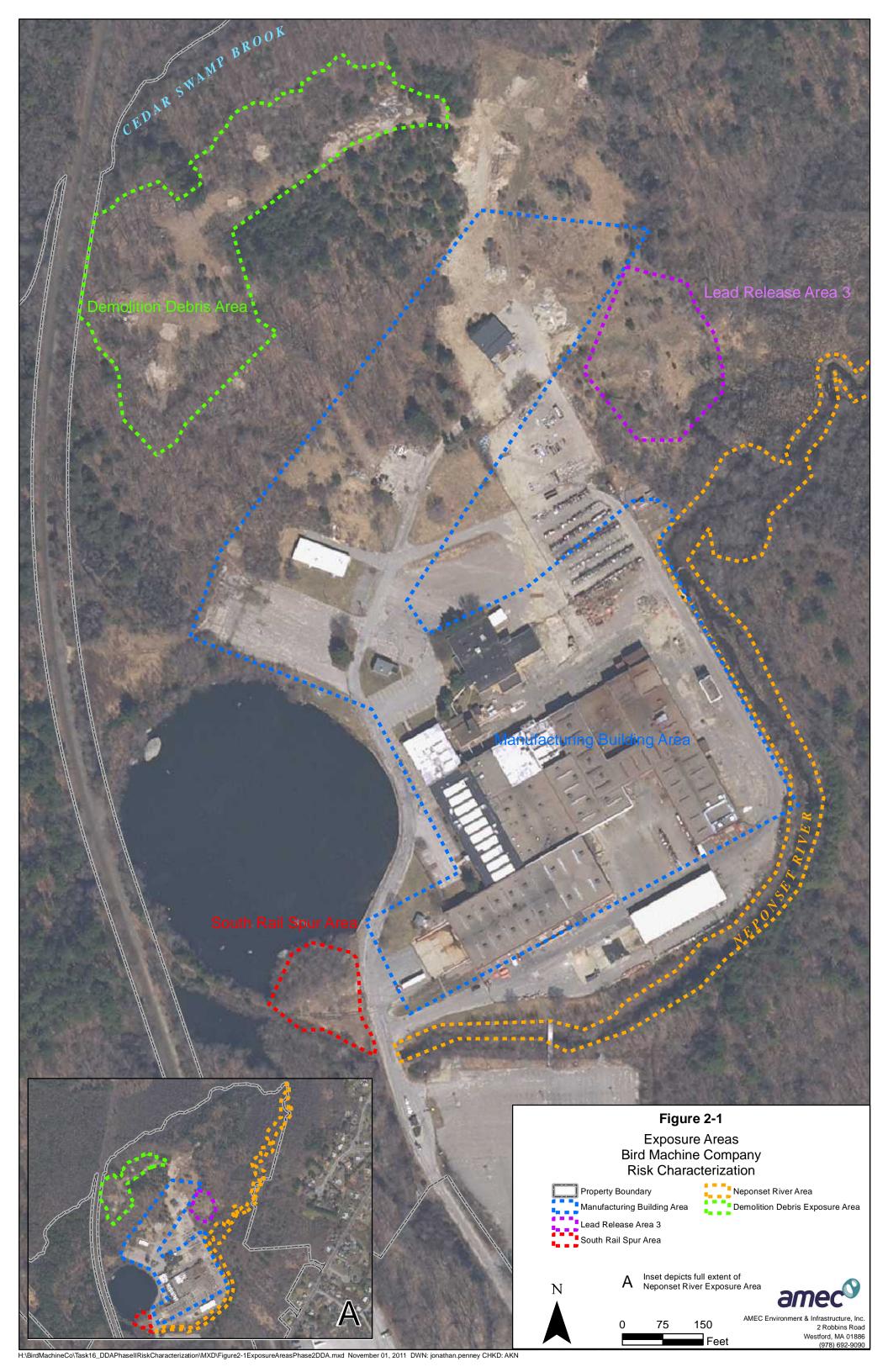
SITE LOCATION MAP

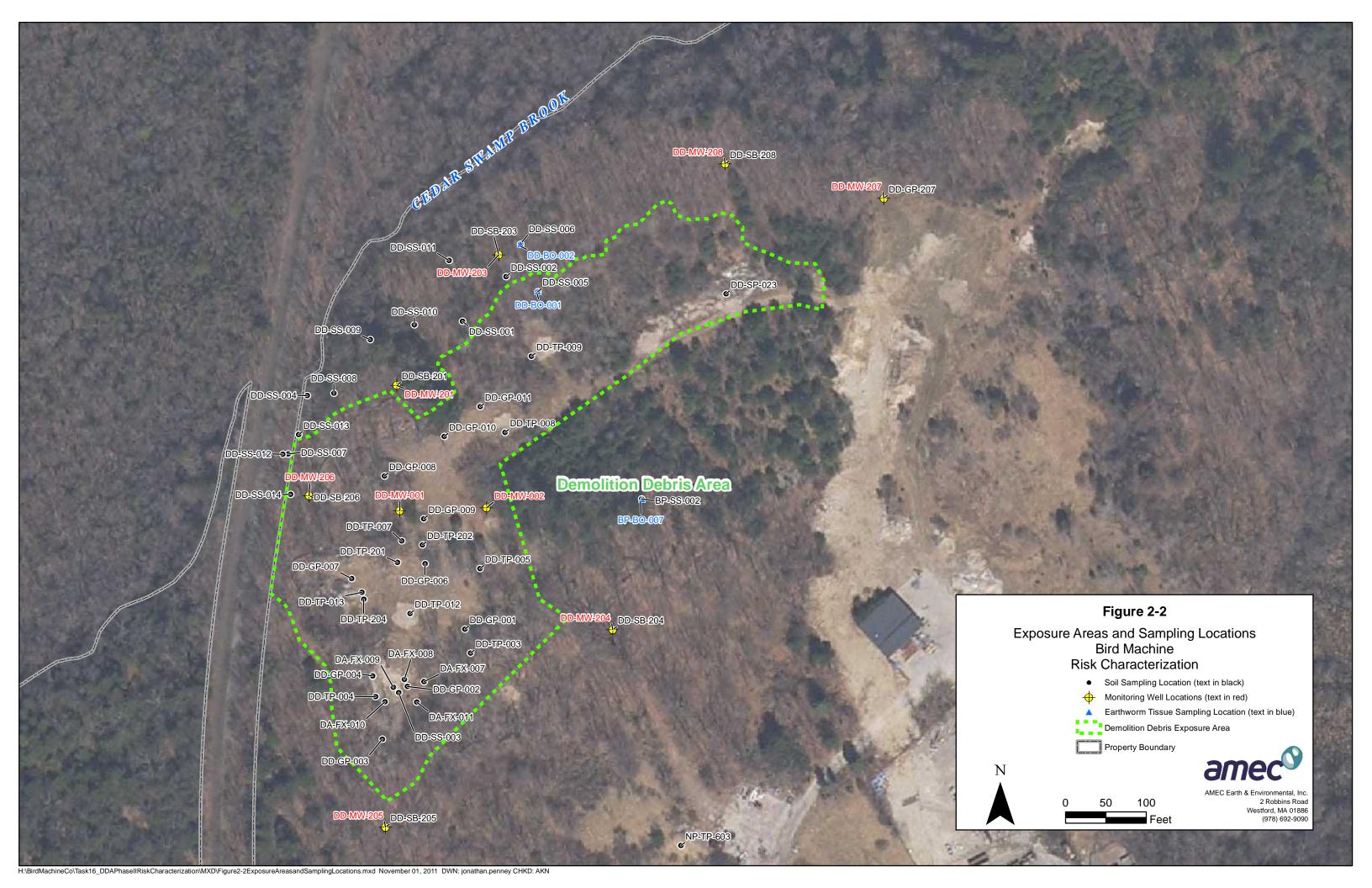
Bird Machine Inc. Company

100 Neponset Street Walpole, MA









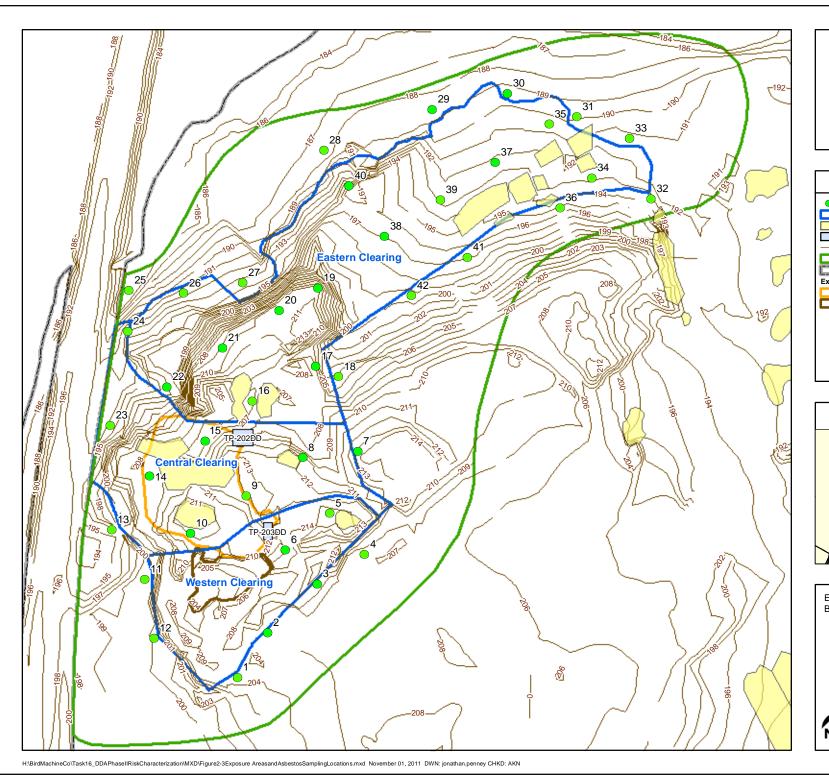


FIGURE 2-3

Asbestos In Soil Sampling Locations

Bird Machine Co. Walpole, MA





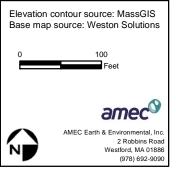
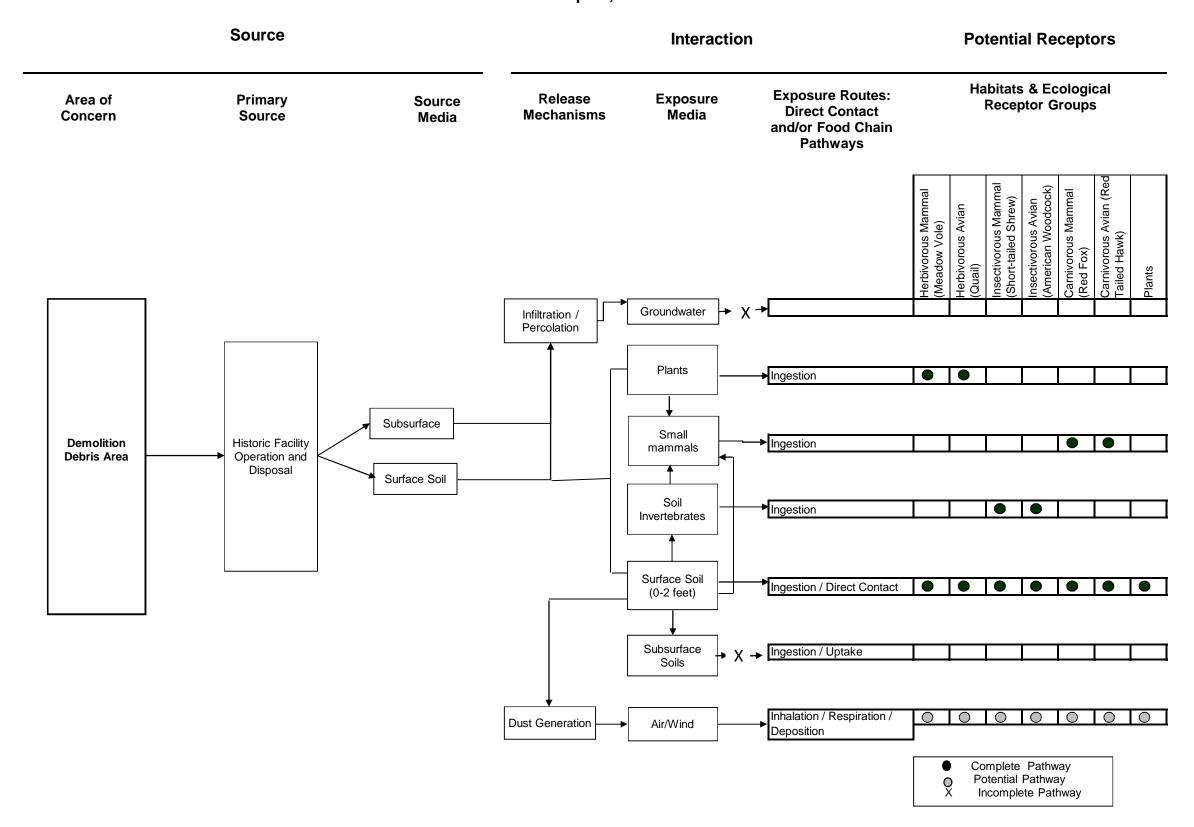


Figure 4-1
Conceptual Site Model
for Environmental Risk Characterization
Bird Machine Company
Walpole, MA



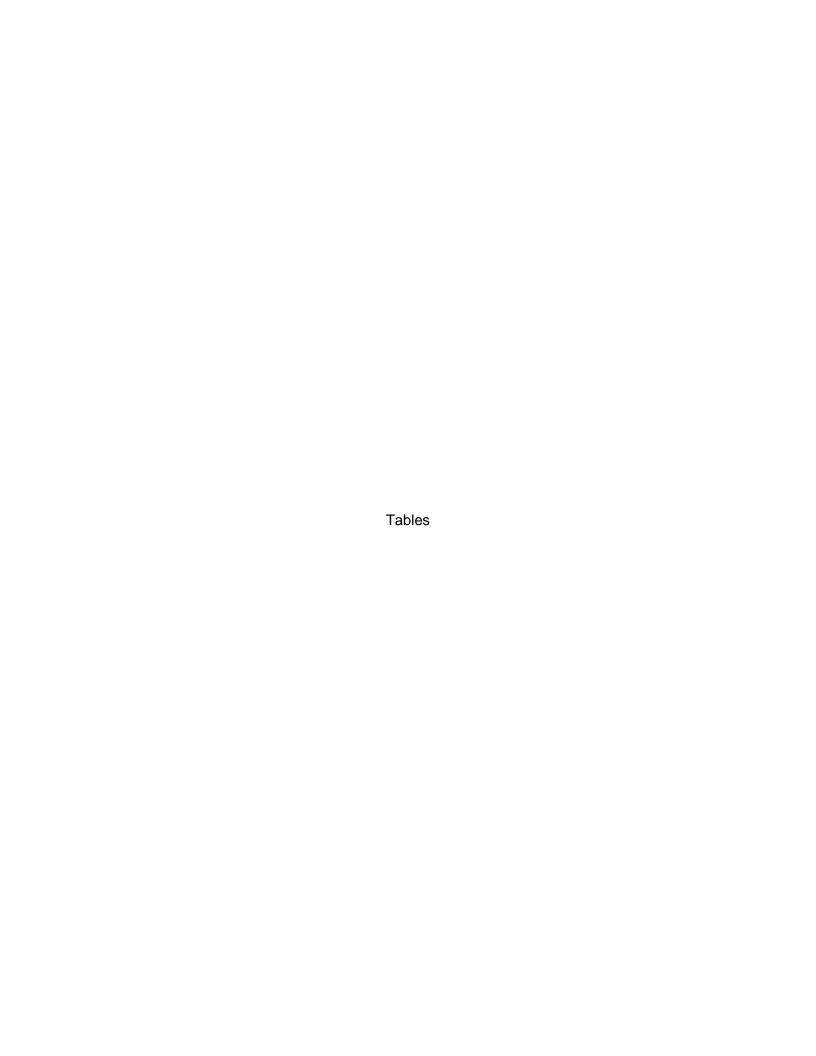


Table 2-1
Soil Samples Included in Exposure Areas
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

Sample ID Number	Sample Used in Ecological Risk Calculations (0-2 ft bgs)	Sample Used in Human Health Risk Calculations (0-3 ft bgs)
DDA-GP-3 (0-2)RR & Dup	X	X
DDA-GP-4 (4-6)	X	X
DDA-GP-6 (0-2)	X	X
DDA-GP-9(4-6)RR & Dup	X	X
DD-GP-207-001-X	,	X
DD-SB-206-003-X	X	X
DD-SS-001-001-X	X	X
DD-SS-005-001-X	X	X
DD-SS-007-001-X & Dup	X	Х
DD-SS-012-001-X	X	Х
DD-SS-014-001-X	X	Х
DD-TP-001-001-X	X	X
DD-TP-001-002-X	X	X
DD-TP03-2S	X	X
DD-TP04-2S	X	X
DD-TP05-2S	X	X
DD-TP07-2S-D & Dup	X	X
DD-TP08-2S	X	X
DD-TP09-2S-D & Dup	X	X
DD-TP12-5W	X	X
DD-TP-201-001-X	X	X
DD-TP-202-002-X	X	X
DD-TP-202-003-X	X	X
SF-TP01-3S		X

Table 3-1
Selection of COCs and Exposure Point Concentrations in Soil
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

CAS		Number	Number	Frequency Of	Minimum	Maximum	Average					
Number	Analyte (mg/kg)	Analyzed	Detected	Detection	Detected	Detected	Concentration		UCL	COC	Rationale	EPC
95-63-6	1,2,4-TRIMETHYLBENZENE	7	1	14%	2.00	2.00	0.29	NA	1000	YES	AB	0.29
107-06-2	1,2-DICHLOROETHANE	7	1	14%	0.0020	0.0020	0.021	NA	6000	YES	AB	0.021
108-67-8	1,3,5-TRIMETHYLBENZENE	6	1	17%	0.68	0.68	0.11	NA	1000	YES	AB	0.11
51-28-5	2,4-DINITROPHENOL	15	0	0%	ND	ND	0.35	NA	900	NO	ND	
78-93-3	2-BUTANONE	7	0	0%	ND	ND	0.17	NA	10000	NO	ND	
91-57-6	2-METHYLNAPHTHALENE	17	2	12%	0.066	0.22	0.14	0.5	5000	NO	BB	
106-47-8	4-CHLOROANILINE	16	2	13%	0.11	0.13	0.64	NA	3000	YES	AB	0.64
99-87-6	4-ISOPROPYLTOLUENE	7	1	14%	0.040	0.040	0.026	NA	1000	YES	AB	0.026
100-02-7	4-NITROPHENOL	16	0	0%	ND	ND	1.68	NA	1000	NO	ND	
83-32-9	ACENAPHTHENE	17	1	6%	0.46	0.46	0.15	0.5	10000	NO	BB	
67-64-1	ACETONE	7	1	14%	0.24	0.24	2.051	NA	10000	YES	AB	2.051
120-12-7	ANTHRACENE	17	3	18%	0.22	1.035	0.22	1.0	10000	NO	BB	
7440-36-0	ANTIMONY	14	5	36%	1.30	8.10	2.18	1.0	300	YES	AB	2.18
	AROCLOR-1260	17	1	6%	1.35	1.35	0.13	NA	100	YES	AB	0.13
7440-38-2	ARSENIC	17	14	82%	1.30	26.00	4.73	20	200	YES	AB	4.73
7440-39-3	BARIUM	17	17	100%	1.70	1400.00	248.57	50	10000	YES	AB	248.57
71-43-2	BENZENE	7	0	0%	ND	ND	0.020	NA	9000	NO	ND	
56-55-3	BENZO(A)ANTHRACENE	17	5	29%	0.13	1.59	0.26	2.0	3000	NO	BB	
50-32-8	BENZO(A)PYRENE	17	4	24%	0.31	1.44	0.26	2.0	300	NO	BB	
205-99-2	BENZO(B)FLUORANTHENE	17	4	24%	0.14	2.30	0.29	2.0	3000	YES	AB	0.29
191-24-2	BENZO(G,H,I)PERYLENE	17	1	6%	1.30	1.30	0.20	1.0	10000	YES	AB	0.20
207-08-9	BENZO(K)FLUORANTHENE	17	4	24%	0.31	2.10	0.30	1.0	10000	YES	AB	0.30
7440-41-7	BERYLLIUM	17	3	18%	0.55	4.10	0.54	0.4	2000	YES	AB	0.54
EPH1122	C11-C22 AROMATICS, ADJUSTED	15	13	87%	11.50	69.00	27.82	NA	1000	YES	AB	27.82
EPH1936	C19-C36 ALIPHATICS	15	13	87%	5.30	170.00	48.070	NA	1000	YES	AB	48.070
VPH58	C5-C8 ALIPHATICS, ADJUSTED	6	1	17%	5.20	5.20	2.042	NA	1000	YES	AB	2.042
VPH910	C9-C10 AROMATICS	6	1	17%	24.00	24.00	5.17	NA	1000	YES	AB	5.17
VPH912	C9-C12 ALIPHATICS, ADJUSTED	6	0	0%	ND	ND	1.41	NA	1000	NO	ND	
VPH918	C9-C18 ALIPHATICS	16	4	25%	4.80	19.00	4.041	NA	1000	YES	AB	4.041
7440-43-9	CADMIUM	17	13	76%	0.38	2.60	0.93	2	300	YES	AB	0.93
7440-47-3	CHROMIUM	17	17	100%	1.20	2200.00	189.49	30	2000	YES	AB	189.49
18540-29-9	CHROMIUM VI	1	0	0%	ND	ND	0.23	30	2000	NO	ND	
218-01-9	CHRYSENE	17	6	35%	0.14	1.69	0.31	2	10000	NO	BB	
7440-50-8	COPPER	1	1	100%	56.00	56.00	56.00	40	1000	YES	AB	56
53-70-3	DIBENZ(A,H)ANTHRACENE	17	1	6%	1.30	1.30	0.20	0.5	1000	YES	AB	0.20
100-41-4	ETHYLBENZENE	7	1	14%	0.23	0.23	0.033	NA	10000	YES	AB	0.033
206-44-0	FLUORANTHENE	17	8	47%	0.16	4.40	0.63	4	10000	YES	AB	0.63
86-73-7	FLUORENE	17	1	6%	0.74	0.74	0.17	1	10000	NO	BB	
193-39-5	INDENO(1,2,3-CD)PYRENE	17	1	6%	0.94	0.94	0.20	1	3000	NO	BB	
7439-92-1	LEAD	17	17	100%	3.00	450.00	100.65	100	3000	YES	AB	101
1330-20-7	M,P-XYLENES	17	2	12%	0.27	0.27	0.18	NA	10000	YES	AB	0.18
7439-97-6	MERCURY	16	10	63%	0.062	2.90	0.42	0.3	300	YES	AB	0.42
91-20-3	NAPHTHALENE	17	3	18%	0.24	0.29	0.11	0.5	10000	NO	BB	
104-51-8	N-BUTYLBENZENE	6	1	17%	0.13	0.13	0.023	NA	1000	YES	AB	0.023
7440-02-0	NICKEL	15	15	100%	0.71	1400.00	224.90	20	7000	YES	AB	225
103-65-1	N-PROPYLBENZENE	7	1	14%	0.28	0.28	0.041	NA	1000	YES	AB	0.041
95-47-6	O-XYLENE	7	1	14%	0.49	0.49	0.070	NA	10000	YES	AB	0.070
85-01-8	PHENANTHRENE	17	7	41%	0.17	4.35	0.53	3	10000	YES	AB	0.53
129-00-0	PYRENE	17	8	47%	0.21	4.050	0.60	4	10000	YES	AB	0.60
7782-49-2	SELENIUM	16	0	0%	ND	ND	0.94	0.5	8000	NO	ND	
7440-22-4	SILVER	17	3	18%	0.83	3.20	1.16	0.6	2000	YES	AB	1.2
98-06-6	TERT-BUTYLBENZENE	7	1	14%	0.24	0.24	0.035	NA	1000	YES	AB	0.035
108-88-3	TOLUENE	7	2	29%	0.14	0.14	0.040	NA	10000	YES	AB	0.040
	ı											

CAS		Number	Number	Frequency Of	Minimum	Maximum	Average					
Number	Analyte (mg/kg)	Analyzed	Detected	Detection	Detected	Detected	Concentration	Background	UCL	COC	Rationale	EPC
7440-62-2	VANADIUM	17	17	100%	1.30	150.00	38.27	30	10000	YES	AB	38
7440-66-6	ZINC	16	16	100%	4.10	810.00	169.19	100	10000	YES	AB	169
1746-01-6	2,3,7,8-TCDD TEQ	12	12	100%	0.00000072	0.000093	0.000027	2.00E-05	0.003	YES	AB	0.000027
NA	Asbestos in Soil (MFG)											
	TOTAL ASBESTOS FIBERS > 5µM	3	3	100%	11.7	82.9	36.6	NA	NA	YES	AB	83

Notes:

COCs - Contaminants of concern

EPC - Exposure Point Concentration

DDA - Drum Disposal Area

RL - Reporting limit

UCL - Upper concentration limit

NA - not applicable/not evaluated

ND - not detected in any sample

AB - above background

BB - below background

mg/kg - milligrams per kilogram

MFG - million fibers per gram

μM - micrometers

(1) Average concentration calculated using 1/2 RLs for non-detects.

For asbestos, the maximum concentration was used as the EPC due to the small data set and high uncertainty associated with predicting airborne concentrations.

(2) Background concentrations obtained from: Background Levels of

Polycyclic Aromatic Hydrocarbons and Metals in Soil (MassDEP, 2002).

- (3) UCLs are the Upper Concentration Limit for soil published at 310 CMR 40.0996(7) Table 6, dated December 2007.
- (4) EPCs are the average concentrations for each analyte.
- (5) The UCLs for total PCBs and total xylenes were used for Aroclor-1260 and m,p-xylenes and o-xylenes, respectively.

Table 3-2 Selection of COCs and Exposure Point Concentrations in Groundwater Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

					All Data			2008 Only										
CAS		Number	Number	Frequency Of	Minimum	Maximum	Average	Number	Number	Frequency Of	Minimum	Maximum	Average					
Number	Analyte (ug/L)	Analyzed	Detected	Detection	Detected	Detected	Concentration	Analyzed	Detected	Detection	Detected	Detected	Concentration	MCL	UCL	COC	Rationale	EPC
51-28-5	2,4-DINITROPHENOL	38	1	3%	0.84	0.84	3.21	9	1	11%	0.84	0.84	2.37		100000	NO	FOD	
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	38	15	39%	0.43	5.3	3.57	9	6	67%	0.43	1.6	1.29	6	100000	YES	AB	1.29
84-74-2	DI-N-BUTYL PHTHALATE	38	1	3%	1.5	1.5	3.26	9	1	11%	1.5	1.5	2.44		10000	NO	FOD	
7440-38-2	ARSENIC (DISSOLVED)	40	10	25%	0.81	10	2.99	9	1	11%	3.9	3.9	0.93	10	9000	YES	AB	0.93
	BARIUM (DISSOLVED)	40	40	100%	6.1	110	30.53	9	9	100%	7.6	46	25.18	2000	100000	YES	AB	25.18
7440-47-3	CHROMIUM (DISSOLVED)	40	24	60%	0.36	43	7.62	9	8	89%	0.36	3.5	1.26	100	3000	YES	AB	1.26
7439-92-1	LEAD (DISSOLVED)	40	1	3%	0.46	0.46	1.22	9	1	11%	0.46	0.46	0.50	15	150	NO	FOD	1
	NICKEL (DISSOLVED)	40	35	88%	0.43	110	5.55	9	9	100%	0.43	2	1.09		2000	YES	AB	1.09
	SELENIUM (DISSOLVED)	40	7	18%	5.5	9.3	3.02	9	2	22%	6.8	7.8	2.07	50	1000	YES	AB	2.07
7440-62-2	VANADIUM (DISSOLVED)	40	8	20%	0.18	120	5.37	9	4	44%	0.18	0.52	0.42		40000	YES	AB	0.42
7440-66-6	ZINC (DISSOLVED)	40	28	70%	1.2	850	28.94	9	8	89%	1.2	110	14.38		50000	YES	AB	14.38

Notes:

MCL - Maximum contaminant level
UCL - Upper concentration limit
COC - Contaminants of concern
EPC - Exposure point concentration
NA - Not applicable/not evaluated
ND - not detected in any sample
AB - above background
BB - below background
FOD - Low frequency of detection
(1) Average concentration calculated using 1/2 RLs for non-detects.
(2) MCLs were obtained from: 2006 Edition of the Drinking Water Standards and Health Advisories (USEPA, 2006).
(3) UCLs are the Upper Concentration Limit for soil published at 310 CMR 40.0996(7) Table 6, dated December 2007.
(4) EPCs are the average 2008 concentrations for each analyte.

Table 3-3
Toxicity Information for Method 3 Human Health Risk Characterization
Bird Machine Company - Walpole, MA
Human Health Risk and Environmental Risk Characterization

	Reference	Reference	Reference	Reference		
	Dose Oral	Dose Oral	Concentration	Concentration	Cancer Slope	Unit Risk
	Chronic	Subchronic	Inhalation	Inhalation	Factor Oral	Inhalation
	Value	Value	Chronic Value	Subchronic	Value (mg/kg-	Value
Analyte	(mg/kg-day)	(mg/kg-day)	(mg/m³)	Value (mg/m³)	day) ⁻¹	(ug/m3) ⁻¹
1,2,4-TRIMETHYLBENZENE	NA	NA	0.007	0.007	NA	NA
1,2-DICHLOROETHANE	0.02	0.2	0.055	0.055	0.091	0.000026
1,3,5-TRIMETHYLBENZENE	0.01	0.01	NA	NA	NA	NA
4-CHLOROANILINE	0.004	0.004	0.014	0.014	0.2	NA
4-ISOPROPYLTOLUENE	NA	NA	NA	NA	NA	NA
ACETONE	0.9	2.7	0.8	0.8	NA	NA
ANTIMONY	0.0004	0.0004	0.01	0.01	NA	NA
AROCLOR-1260 (TOTAL PCBs)	0.00002	0.00005	0.00002	0.00002	2	1.0E-04
ARSENIC	0.0003	0.0003	0.0000025	0.0000025	1.5	0.0043
BARIUM	0.2	0.07	0.0005	0.005	NA	NA
BENZO(B)FLUORANTHENE	0.03	0.3	0.05	0.5	0.73	0.00011
BENZO(G,H,I)PERYLENE	0.03	0.3	0.05	0.5	NA	NA
BENZO(K)FLUORANTHENE	0.03	0.3	0.05	0.5	0.073	0.00011
BERYLLIUM	0.002	0.005	0.00002	0.00002	NA	0.0024
C11-C22 AROMATICS	0.03	0.3	0.05	0.5	NA	NA
C19-C36 ALIPHATICS	2	6	NA	NA	NA	NA
C5-C8 ALIPHATICS	0.04	0.4	0.2	0.2	NA	NA
C9-C10 AROMATICS	0.03	0.3	0.05	0.5	NA	NA
C9-C18 ALIPHATICS	0.1	1	0.2	0.6	NA	NA
CADMIUM	0.0005	0.0005	0.00002	0.00002	NA	0.00180
CHROMIUM (III)	1.5	1.5	0.0001	0.0003	NA	NA
COPPER	0.04	NA	NA	NA	NA	NA
DIBENZ(A,H)ANTHRACENE	0.03	0.3	0.05	0.5	7.3	0.0012
ETHYLBENZENE	0.1	1	1	1	0.11	0.0000025
FLUORANTHENE	0.04	0.4	0.05	0.5	NA	NA
LEAD	0.00075	0.00075	0.001	0.001	NA	NA
M,P-XYLENES (TOTAL XYLENES)	0.2	0.2	0.1	0.3	NA	NA
MERCURY	0.0003	0.0003	0.0003	0.0003	NA	NA
N-BUTYLBENZENE	0.05	0.05	NA	NA	NA	NA
NICKEL	0.02	0.02	0.001	0.001	NA	0.00026
N-PROPYLBENZENE	0.1	0.1	1	1	NA	NA
O-XYLENE (TOTAL XYLENES)	0.2	0.2	0.1	0.3	NA	NA

Table 3-3
Toxicity Information for Method 3 Human Health Risk Characterization
Bird Machine Company - Walpole, MA
Human Health Risk and Environmental Risk Characterization

Analyte	Reference Dose Oral Chronic Value (mg/kg-day)	Reference Dose Oral Subchronic Value (mg/kg-day)	Concentration Inhalation Chronic Value (mg/m³)	Concentration Inhalation Subchronic Value (mg/m³)	Cancer Slope Factor Oral Value (mg/kg- day) ⁻¹	Unit Risk Inhalation Value (ug/m3) ⁻¹
PHENANTHRENE	0.03	0.3	0.05	0.5	NA	NA
PYRENE	0.03	0.3	0.05	0.5	NA	NA
SILVER	0.005	0.005	0.00014	0.00014	NA	NA
TERT-BUTYLBENZENE	NA	NA	NA	NA	NA	NA
TOLUENE	0.08	0.8	5	5	NA	NA
VANADIUM	0.009	0.001	0.001	0.001	NA	NA
ZINC	0.3	0.3	0.0014	0.0014	NA	NA
2,3,7,8-TCDD TEQ	1E-09	1E-09	0.00000004	0.00000004	130000	38

	Reference	Reference	Reference	Reference		
	Dose Oral	Dose Oral	Concentration	Concentration	Cancer Slope	Unit Risk
	Chronic	Subchronic	Inhalation	Inhalation	Factor Oral	Inhalation
	Value	Value	Chronic Value	Subchronic	Value (mg/kg-	Value
Analyte	(mg/kg-day)	(mg/kg-day)	(mg/m³)	Value (mg/m ³)	day) ⁻¹	(fibers/ml) ⁻¹
ASBESTOS	NA	NA	NA	NA	NA	0.23

mg/kg-day - milligrams per kilogram per day

mg/m3 - milligram per cubic meter

(ug/m3)-1 - inverse of microgram per cubic meter

NA - not available

- (1) All un-shaded cells contain toxicity data from the December 2009 MassDEP Toxicity Spreadsheet.
- (2) Gray cells contain toxicity data from USEPA (June 2011 RSLs for chemicals and the IRIS profile for asbestos).
- (3) Benzo(b)fluoranthene, benzo(g,h,i)perylene and dibenzo(a,h)anthracene have calculated MassDEP IURs but RSLs were used in lieu of the calculated values.
- (4) The values for total PCBs were used for Aroclor-1260.
- (5) The value for chromium III were used for chromium.
- (6) The values for total xylenes were used for m,p-xylenes and o-xylenes.

Table 3-4
World Health Organization Toxicity Equivalency Factors for Dioxin/Furan Congeners
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Mammalian TEF	Avian TEF
Chlorinated dibenzo-p-dioxins		
2,3,7,8-TCDD	1	1
1,2,3,7,8-PeCDD	1	1
1,2,3,4,7,8-HxCDD	0.1	0.05
1,2,3,6,7,8-HxCDD	0.1	0.01
1,2,3,7,8,9-HxCDD	0.1	0.01
1,2,3,4,6,7,8-HpCDD	0.01	0.001
OCDD	0.0003	0.0001
Chlorinated dibenzofurans		
2,3,7,8-TCDF	0.1	1
1,2,3,7,8-PeCDF	0.03	0.01
2,3,4,7,8-PeCDF	0.3	1
1,2,3,4,7,8-HxCDF	0.1	0.1
1,2,3,6,7,8-HxCDF	0.1	0.1
1,2,3,7,8,9-HxCDF	0.1	0.1
2,3,4,6,7,8-HxCDF	0.1	0.1
1,2,3,4,6,7,8-HpCDF	0.01	0.01
1,2,3,4,7,8,9-HpCDF	0.01	0.01
OCDF	0.0003	0.0001

TEF - Toxicity equivalency factor

TEFs presented are the World Health Organization (WHO) 2005 values from Van den Berg et al., 2006

Table 3-5 Relative Absorption Factors
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Relative	Relative	Relative	Relative				Relative	Relative
	Absorption	Absorption	Absorption	Absorption	Relative		Relative	Absorption	Absorption
	Factor Soil	Factor Soil	Factor Soil	Factor Soil	Absorption	Relative	Absorption	Factor Air	Factor Air
	Ingestion Non-	Dermal Non-	Ingestion Non-	Dermal Non-	Factor Soil	Absorption	Factor Air		Inhalation Non-
	Cancer	Cancer	Cancer	Cancer	Ingestion	Factor Soil	Inhalation	Cancer	Cancer
Compound	Chronic	Chronic	Subchronic	Subchronic	Cancer	Dermal Cancer	Cancer	Chronic	Subchronic
1,2,4-TRIMETHYLBENZENE	1	1	1	1	NC	NC	NC	1	1
1,2-DICHLOROETHANE	1	0.1	1	0.1	1	0.1	1	1	1
1,3,5-TRIMETHYLBENZENE	1	1	1	1	NC	NC	NC	1	1
4-CHLOROANILINE	1	0.1	1	0.1	1	0.08	1	1	1
4-ISOPROPYLTOLUENE	1	0.12	1	0.12	NC	NC	NC	1	1
ACETONE	1	0.1	1	0.1	NC	NC	NC	1	1
ANTIMONY	1	0.1	1	0.1	NC	NC	NC	1	1
AROCLOR-1260	0.85	0.16	0.85	0.16	0.85	0.16	1	1	1
ARSENIC	1	0.03	1	0.03	1	0.03	1	1	1
BARIUM	1	0.05	1	0.05	NC	NC	NC	1	1
BENZO(B)FLUORANTHENE	0.28	0.02	0.28	0.02	0.28	0.02	1	1	1
BENZO(G,H,I)PERYLENE	0.36	0.1	0.36	0.1	NC	NC	NC	1	1
BENZO(K)FLUORANTHENE	0.28	0.02	0.28	0.02	0.28	0.02	1	1	1
BERYLLIUM	1	0.03	1	0.03	NC	0.03	1	1	1
C11-C22 AROMATICS	0.36	0.1	0.36	0.1	NC	NC	NC	1	1
C19-C36 ALIPHATICS	1	0.1	1	0.1	NC	NC	NC	1	1
C5-C8 ALIPHATICS	1	1	1	1	NC	NC	NC	1	1
C9-C10 AROMATICS	1	0.5	1	0.5	NC	NC	NC	1	1
C9-C18 ALIPHATICS	1	0.5	1	0.5	NC	NC	NC	1	1
CADMIUM	1	0.14	1	0.14	NC	NC	1	1	1
CHROMIUM	1	0.04	1	0.04	NC	NC	NC	1	1
COPPER	1	1	1	1	NC	NC	1	1	1
DIBENZ(A,H)ANTHRACENE	0.28	0.02	0.28	0.02	0.28	0.02	1	1	1
ETHYLBENZENE	1	0.2	1	0.2	1	0.08	1	1	1
FLUORANTHENE	0.36	0.1	0.36	0.1	NC	NC	NC	1	1
LEAD	0.5	0.006	0.5	0.006	NC	NC	NC	1	1
M,P-XYLENES	1	0.12	1	0.12	NC	NC	NC	1	1
MERCURY	1	0.05	1	0.05	NC	NC	NC	1	1
N-BUTYLBENZENE	1	1	1	1	NC	NC	NC	1	1
NICKEL	1	0.35	1	0.35	NC	NC	1	1	1
N-PROPYLBENZENE	1	1	1	1	NC	NC	NC	1	1
O-XYLENE	1	0.12	1	0.12	NC	NC	NC	1	1
PHENANTHRENE	0.36	0.1	0.36	0.1	NC	NC	NC	1	1
PYRENE	0.36	0.1	0.36	0.1	NC	NC	NC	1	1
SILVER	1	0.25	1	0.25	NC	NC	NC	1	1
TERT-BUTYLBENZENE	1	1	1	1	NC	NC	NC	1	1
TOLUENE	1	0.12	1	0.12	NC	NC	NC	1	1
VANADIUM	1	0.03	1	0.03	NC	NC	NC	1	1
ZINC	1	0.02	1	0.02	NC	NC	NC	1	1
2,3,7,8-TCDD TEQ	1	1	1	1	1	0.2	1	1	1
ASBESTOS	-	-	-	•	-	-	1	-	-

Notes:
NC - Not calculated
The Soil Cancer Ingestion, Soil Cancer Dermal, and Inhalation Cancer RAFs for benzene were used as surrogates for p-chloroaniline and ethylbenzene.

Table 3-6
Exposure Assumptions
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

				Trespasser Older Child		
				Older Cillia		
Medium	Route	Parameter	Units	Value	Source	Comment
	Incidental Ingestion &					
Soil	Dermal Contact	Soil Ingestion Rate	mg/d	50	С	
		Soil TWA Ingestion Rate	mg-y/kg-d			
						Median values for face, hands, forearms, lower legs, and feet of females; average from age 8
		Soil Dermal Contact Skin Exposed	cm2/d	4260	а	to 15.
		Soil TWA Dermal Contact Skin Exposed	mg-cm2-y/kg-d-cm2			
		Soil Dermal Contact Adherence Rate	mg/cm2	0.14	С	
		Soil Exposure Frequency	d/y	50	b	
		Soil Exposure Period - Cancer	у	7	С	
		Soil Exposure Period - Non-Cancer	у	7	С	
		Soil Averaging Time - Cancer	d	25550	С	
		Soil Averaging Time - Non-Cancer	d	2555	С	
General		Body Weight	kg	39.9	С	
			_			

mg/d - milligrams per day

mg-y/kg-d - milligram per year per kilogram per day

cm2/d - square centimeters per day

y - year

d - day

kg - kilogram

a - MADEP (1995a). Appendices to the Guidance for Disposal Site Risk Characterization - In support of the Massachusetts Contingency Plan. Interim Final Policy. WSC/ORS-95-141. July, 1995.

b - Professional judgement

c - Value for resident

$$CR_{\text{TWA}} = \frac{CR_{\text{YC}} \times EP_{\text{YC}}}{BW_{\text{YC}}} + \frac{CR_{\text{OC}} \times EP_{\text{OC}}}{BW_{\text{OC}}} + \frac{CR_{\text{A}} \times EP_{\text{A}}}{BW_{\text{A}}}$$

where :

CR = Contact Rate

EP = Exposure Period

BW = Body Weight

TWA = Time-Weighted Average

YC = Young Child

OC = Older Child

A = Adult

Table 3-7
Estimated Potential Human Health Risks - Soil
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Non-Cancer	Cancer	Asbestos
Receptor	HI	Risk	Risk
Current Trespasser (0-3') DDA	0.10	9E-06	8.4E-06
Current Trespasser (0-3') SFA	0.03	5E-10	NC

HI - Hazard Index DDA - Demolition Debris Area SFA - Sand Filter Area NC - not calculated

Table 3-8
Estimated Potential Human Health Risks - Groundwater
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

Summary of Total HQs

COC	DD-MW-002	DD-MW-201	DD-MW-203	DD-MW-204	DD-MW-205	DD-MW-206	DD-MW-207
Bis(2-ethylhexyl)phthalate	0.0092	0.0056	0.0017	0.0018	0.0092	0.0020	0.0020
Barium	0.0056	0.013	0.0033	0.0059	0.0023	0.0047	0.0092
Chromium (total)	0.041	0.059	0.016	0.013	0.014	0.014	0.029
Nickel	0.0036	0.0037	0.0036	0.0036	0.0018	0.0021	0.0013
Selenium	0.0059	0.086	0.0059	0.012	0.0059	0.0059	0.0059
Vanadium	0.0036	0.0026	0.0020	0.0036	0.0023	0.0036	0.0036
Zinc	0.00024	0.00049	0.0010	0.00049	0.00032	0.00037	0.00037
Total	0.07	0.17	0.03	0.04	0.04	0.03	0.05

Summary of Total ELCRs

COC(1)	DD-MW-002	DD-MW-201	DD-MW-203	DD-MW-204	DD-MW-205	DD-MW-206	DD-MW-207
Bis(2-ethylhexyl)phthalate	8E-07	5E-07	1E-07	1E-07	8E-07	2E-07	2E-07
Total	8E-07	5E-07	1E-07	1E-07	8E-07	2E-07	2E-07

Notes:

COC - contaminant of concern

HQ - Hazard quotient

ELCR - Excess lifetime cancer risk

1. All other COCs are not considered to be carcinogenic

Table 4-1
Exposure Point Concentrations and Selection of COPECs in DDA Soil
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

		Number	Number	Detection	Minimum	Maximum	Average	MCP	Lowest	Rationale	000500	EDO (4)
CAS Number	Analyte	Analyzed	Detected	Frequency	Detected	Detected	Concentration (1)	Background (2)	Eco SSL	(3)	COPEC?	EPC (4)
107-06-2	1,2-DICHLOROETHANE	7	1	14.29%	0.0020	0.0020	0.021	NA ()	21.2	BB	no	NA
78-93-3	2-BUTANONE	7	0	0.00%	ND	ND	0.17	NA	89.6	ND	no	NA
91-57-6	2-METHYLNAPHTHALENE	16	2	12.50%	0.066	0.22	0.14	NA	29	BB	no	NA
106-47-8	4-CHLOROANILINE	15	2	13.33%	0.11	0.13	0.65	NA	1.1	BB	no	NA
83-32-9	ACENAPHTHENE	16	1	6.25%	0.46	0.46	0.16	0.5	20	BB	no	NA
67-64-1	ACETONE	7	1	14.29%	0.24	0.24	2.051	NA	2.5	BB	no	NA
120-12-7	ANTHRACENE	16	3	18.75%	0.22	1.035	0.23	1	29	BB	no	NA
7440-36-0	ANTIMONY	13	5	38.46%	1.30	8.10	2.24	1	0.27	AB, ASSL	yes	2.24
11096-82-5	AROCLOR-1260	16	1	6.25%	1.35	1.35	0.14	NA	0.000332	ASSL	ves	0.14
7440-38-2	ARSENIC	16	13	81.25%	1.30	26.00	4.84	20	18	AB, ASSL	yes	4.84
7440-39-3	BARIUM	16	16	100.00%	1.70	1400.00	262.42	50	330	AB, ASSL	yes	262.42
71-43-2	BENZENE	7	0	0.00%	ND	ND	0.020	NA	0.255	ND	no	NA
56-55-3	BENZO(A)ANTHRACENE	16	5	31.25%	0.13	1.59	0.27	2	1.1	BB	no	NA
50-32-8	BENZO(A)PYRENE	16	4	25.00%	0.31	1.44	0.28	2	1.1	BB	no	NA
205-99-2	BENZO(B)FLUORANTHENE	16	4	25.00%	0.14	2.30	0.30	2	1.1	AB, ASSL	ves	0.30
191-24-2	BENZO(G,H,I)PERYLENE	16	1	6.25%	1.30	1.30	0.21	1	1.1	AB, ASSL	ves	0.21
207-08-9	BENZO(K)FLUORANTHENE	16	4	25.00%	0.31	2.10	0.31	1	1.1	AB, ASSL	ves	0.31
7440-41-7	BERYLLIÚM	16	3	18.75%	0.55	4.10	0.57	0.4	10	BB	no	NA
EPH1122	C11-C22 AROMATICS, ADJUSTED	14	13	92.86%	11.5	69	29.67	NA	NA	NA	ves	29.67
EPH1936	C19-C36 ALIPHATICS	14	13	92.86%	5.3	170	51.37	NA	NA	NA	yes	51.37
VPH58	C5-C8 ALIPHATICS, ADJUSTED	6	1	16.67%	5.2	5.2	2.042	NA	NA	NA	yes	2.042
VPH910	C9-C10 AROMATICS	6	1	16.67%	24	24	5.17	NA	NA	NA	ves	5.17
VPH912	C9-C12 ALIPHATICS, ADJUSTED	6	0	0.00%	ND	ND	1.41	NA	NA	ND	no	NA
VPH912	C9-C18 ALIPHATICS	15	4	26.67%	4.8	19	4.18	NA	NA	NA	ves	4.18
7440-43-9	CADMIUM	16	13	81.25%	0.38	2.60	0.98	2	0.36	AB, ASSL	ves	0.98
7440-47-3	CHROMIUM	16	16	100.00%	1.20	2200.00	200.27	30	0.4	AB, ASSL	ves	200.27
18540-29-9	CHROMIUM VI	1	0	0.00%	ND	ND	0.23	30	NA	ND	no	NA
218-01-9	CHRYSENE	16	6	37.50%	0.14	1.69	0.33	2	1.1	BB	no	NA
7440-50-8	COPPER	1	1	100.00%	56.00	56.00	56.00	40	28	AB, ASSL	ves	56.00
53-70-3	DIBENZ(A,H)ANTHRACENE	16	1	6.25%	1.30	1.30	0.21	0.5	1.1	AB, ASSL	yes	0.21
100-41-4	ETHYLBENZENE	7	1	14.29%	0.23	0.23	0.033	NA	5.16	BB	no	NA
206-44-0	FLUORANTHENE	16	8	50.00%	0.16	4.40	0.66	4	29	BB	no	NA
86-73-7	FLUORENE	16	1	6.25%	0.74	0.74	0.17	1	29	BB	no	NA
193-39-5	INDENO(1,2,3-CD)PYRENE	16	1	6.25%	0.94	0.94	0.20	1	1.1	BB	no	NA
7439-92-1	LEAD	16	16	100.00%	3.00	450.00	106.47	100	11	AB, ASSL	yes	106.47
0	M,P-XYLENES	16	2	12.50%	0.27	0.27	0.18	NA	10	BB	no	NA
7439-97-6	MERCURY	15	10	66.67%	0.062	2.90	0.44	0.3	0.1	AB	yes	0.44
91-20-3	NAPHTHALENE	16	3	18.75%	0.24	0.29	0.12	0.5	29	BB	no	NA
104-51-8	N-BUTYLBENZENE	6	1	16.67%	0.13	0.13	0.023	NA	5.16	BB	no	NA
7440-02-0	NICKEL	14	14	100.00%	0.71	1400.00	239.82	20	38	AB, ASSL	yes	239.82
103-65-1	N-PROPYLBENZENE	7	1	14.29%	0.28	0.28	0.041	NA	5.16	BB	no	NA
95-47-6	O-XYLENE	7	1	14.29%	0.49	0.49	0.070	NA	10	BB	no	NA
85-01-8	PHENANTHRENE	16	7	43.75%	0.17	4.35	0.56	3	29	BB	no	NA
129-00-0	PYRENE	16	8	50.00%	0.21	4.050	0.63	4	1.1	AB, ASSL	yes	0.63
7782-49-2	SELENIUM	15	0	0.00%	ND	ND	0.96	0.5	0.52	ND	no	NA
7440-22-4	SILVER	16	3	18.75%	0.83	3.20	1.19	0.6	4.2	BB	no	NA

Table 4-1
Exposure Point Concentrations and Selection of COPECs in DDA Soil
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

		Number	Number	Detection	Minimum	Maximum	Average	MCP	Lowest	Rationale	COPEC?	EPC (4)
CAS Number	Analyte	Analyzed	Detected	Frequency	Detected	Detected	Concentration (1)	Background (2)	Eco SSL	(3)	COFEC	EFC (4)
108-88-3	TOLUENE	7	2	28.57%	0.14	0.14	0.040	NA	5.45	BB	no	NA
7440-62-2	VANADIUM	16	16	100.00%	1.30	150.00	38.85	30	2	AB, ASSL	yes	38.85
7440-66-6	ZINC	15	15	100.00%	4.10	810.00	177.61	100	46	AB, ASSL	yes	177.61
1746-01-6	2,3,7,8-TCDD TEQ (mammalian)	12	12	100.00%	0.00000075	0.000095	0.000028	NA	1.99E-07	ASSL	yes	0.000027
1746-01-6	2,3,7,8-TCDD TEQ (avian)	12	12	100.00%	0.00000110	0.00018	0.000049	NA	1.99E-07	ASSL	yes	0.000045

COPEC - Contaminants of potential ecological concern

EPC - Exposure point concentration; represented by average concentrations

DDA - Drum disposal area

NA - not applicable/not evaluated

ND - not detected in any sample AB - above background

BB - below background

ASSL - Above lowest Eco-SSL

MCP - Massachusetts contingency plan

Eco SSL - Ecological soil screening level

- (1) Average concentration was calculated using 1/2 RLs for non-detects.
- (2) Background concentrations were obtained from: Background levels of Polycyclic Aromatic Hydrocarbons and Metals in Soil (MassDEP, 2002).
- (3) Maximum detected concentrations were screened against MCP background as well as lowest Eco SSL for each analyte.

Table 4-3
Potential Exposure Parameter Values for Ecological Receptors
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Avian	Carnivore	Avian	Insectivore	Avian	Herbivore	Mammal	Carnivore	Mammal	Insectivore	Mammal	Herbivore
		ed-Tailed Hawk		ican Woodcock	`	thern Bobwhite)		Fox		ort-tailed Shrew		eadow Vole
Parameter	Value		Value		Value		Value		Value		Value	
Body Weight (kg)	1.134	USEPA (1993), average of six adults	0.178	USEPA (1993) avg of A M & A F in central MA	0.174	USEPA (1993) avg of A M and A F	4.54	USEPA (1993). Mean of male and female BWs in spring and fall.	0.01681	USEPA (1993)	0.0373	USEPA (1993) avg of A M & F all year
Total Dietary Intake (kg ww/d)	0.1191	USEPA (1993) average of A M & F winter	0.1371	USEPA (1993)	0.0134	USEPA (1993)	0.510	USEPA (1993).	0.008	USEPA (1993) A M&F	0.01119	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.00371	Calc. from Beyer (1994)	0.00532	Calc. from Beyer (1994)	0.00017	Calc. from Beyer (1994)	0.00428	Calc. from Beyer (1994)	0.00006	Calc. from Beyer (1994)	0.00003	Calc. from Beyer (1994)
Vegetation Ingestion Rate (kg ww/d)	NR		NR		0.0134	Assume 100% of diet.	NR		NR		0.01119	Assume 100% of diet.
Small Mammal Ingestion Rate (kg ww/d)	0.1191	Assumed 100% of diet	NR		NR		0.50999	Assume 100% of diet.	NR		NR	
Invert Ingestion Rate (kg ww/d)	NR		0.1371	Assumed 100% of diet	NR		NR		0.008	Assumed to be 100% of diet	NR	
Fraction Soil in Diet (kg soil dw/kg diet dw)	0.1040	Beyer et al, (1994) as cited in USEPA (1999) value for woodcock assumed	0.1040	Beyer et al, (1994) as cited in USEPA (1999)	0.1040	Beyer et al, (1994) as cited in USEPA (1999)	0.0280	Beyer et al, (1994)	0.020	Beyer et al, (1994) as cited in USEPA (1999) for deer mouse	0.024	Beyer et al, (1994) as cited in USEPA (1999)
Soil Dry wt./wet wt. CF	0.786	Soils comprised of 78.6% solids.	0.786	Soils comprised of 78.6% solids.	0.786	Soils comprised of 78.6% solids.	0.786	Soils comprised of 78.6% solids.	0.786	Soils comprised of 78.6% solids.	0.786	Soils comprised of 78.6% solids.
Veg Dry wt./wet wt. CF	NR		NR		0.12	USEPA (1999)	NR		NR		0.12	USEPA (1999)
Sm. Mammal Dry wt./wet wt. CF	0.3	USEPA (1999)	NR		NR		0.3	USEPA (1999)	NR		NR	
Invert Dry wt./wet wt. CF	NR		0.373	Site-specific data.	NR		NR		0.373	Site-specific data.	NR	
Home range (ha)	60	USEPA (1993) A M & F spring	3.8	USEPA (1993) avg of inactive A M and brooding A F	9.98	USEPA (1993) avg of A M and A F	1038	USEPA (1993)	0.39	USEPA (1993)	0.06	USEPA (1993) avg of A M & F grassy meadow MA
Area Use Factor (DDA)	0.0308	1.85. ha exposure area	0.4868	1.85. ha exposure area	0.185	1.85. ha exposure area	0.0018	1.85. ha exposure area	1	1.85. ha exposure area	1	1.85. ha exposure area
Area Use Factor (SFA)	0.0308	1.85. ha exposure area	0.00153333	0.092 ha exposure area	0.024210526	0.092 ha exposure area	0.009218437	0.092 ha exposure area	8.9E-05	0.092 ha exposure area	1	0.092 ha exposure area

- (1) USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.
- (2) USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187
- (3) Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382.
- (4) MADEP. 2003. Updated Patroleum Hydrocarbon Fraction Toxicity Values for the VPH/EPH/APH Methodology.

Table 4-4
Soil-to-Biota Transfer Factors
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Invertebra	te	Vegetatio	n	Small Mamn	nal
Constituent	BTFs		BTFs		BTFs	
Antimony	1.00E+00	(g)		(g)		(g)
Arsenic		(g)	3.75E-02	(g)		(g)
Aroclor-1260	1.13E+00	(a)	1.00E-02	(a)	5.83E-05	(e)
Barium	9.10E-02	(g)	1.56E-01	(g)		(g)
Benzo(b)fluoranthene	2.60E+00	(g)	3.10E-01	(g)	0.00E+00	(g)
Benzo(g,h,i)perylene	2.94E+00	(g)		(g)	0.00E+00	(g)
Benzo(k)fluoranthene	2.60E+00	(g)		(g)	0.00E+00	(g)
C11-C22 Aromatics	1.00E+00	(c)	1.00E+00	(c)	1.00E+00	(c)
C19-C36 Aliphatics	1.00E+00	(c)	1.00E+00	(c)	1.00E+00	(c)
C5-C8 Aliphatics	1.00E+00	(c)	1.00E+00	(c)	1.00E+00	(c)
C9-C10 Aromatics	1.00E+00	(c)	1.00E+00	(c)	1.00E+00	(c)
C9-C18 Aliphatics	1.00E+00	(c)	1.00E+00	(c)	1.00E+00	(c)
Cadmium		(g)		(g)		(g)
Chromium	3.06E-01	(g)	4.10E-02	(g)		(g)
Copper	5.15E-01	(g)		(g)		(g)
Dibenz(a,h)anthracene	2.31E+00	(g)	1.30E-01	(g)	0.00E+00	(g)
Lead		(g)		(g)		(g)
Mercury	4.00E-02	(a)	4.30E-02	(a)	7.52E-06	(b)
Nickel	2.70E-02	(a)		(g)		(g)
Pyrene	1.75E+00	(g)	7.20E-01	(g)	0.00E+00	(g)
Vanadium	4.20E-02	(g)	4.85E-03	(g)	1.23E-02	(g)
Zinc		(g)		(g)		(g)
2,3,7,8-TCDD TEQ	1.45E+00	(d)	5.60E-03	(b)	7.81E-05	(b)

References

- (a) USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.
- (b) USEPA (1999). Value for largest BTF in PAH class.
- (c) Conservative default of 1.
- (d) Dioxin insect BTF from Meyn, Ossi, Maurice Zeeman, Michael J Wise, and Susan E. Keane. 1997. Terrestrial Wildlife Risk Assessment for TCDD in Land-Applied Pulp and Paper Mill Sludge. Environmental Toxicology and Chemistry, Vol 16, No. 9, pp 1789-18 (e) USEPA (1999) for deer mouse. Value for Aroclor 1254
- (f) Geomean of BTFs of all other metals.
- (g) USEPA. 2003. Guidance for Developing Ecological Soil Screening Levels. (Eco-SSLs). OSWER Directive 9285.7-55.
- -When BTFs were not available, EcoSSL uptake equations were used. See attachment F-3 for these values
- -All units are in dry weight organism/dry weight soil with the exception of Aroclor 1260, Mercury, Nickel and 2,3,7,8-TCDD. These analytes are in wet weight organism/dry weight soil. Values are adjusted accordingly in risk calculations.

Table 4-5 Mammalian Toxicity Reference Values Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

NOAEL-Based TRVs

NOAEL-Based TRVs	Test Species	Test Species	1				Subchronic	LOAEL	Endpoint	Carnivorous	Herbivorous	Insectivorous	1	1
Chemical	Common Name	Body Weight	Endpoint	Study Type	Effect to Test Organism	Toxicity Value	to Chronic UF	to NOAEL UF	Adjusted TRV	mammal Red Fox (mg/kg-day)	Mammal Meadow Vole (mg/kg-day)	Mammal Short-tailed Shrew (mg/kg-day)	Toxicity Value Surrogate	Initial Compilation Source
		(kg)				(mg/kg-day)			(mg/kg-day)					
	mouse	0.03	NOAEL	chronic	growth, reproduction	13.3			13.3	3.79E+00	1.26E+01	1.54E+01		EPA 2005 (Antimony Eco SSL, geomean of TRVs
Antimony		0.044	NOAEL		1 0	0.000			0.000	4.005.00	5.005.00	0.505.00	A 1 1051	for growth & reproduction)
Aroclor-1260	mouse	0.014	NOAEL	chronic	reproduction	0.068			0.068	1.60E-02	5.32E-02	6.50E-02	Aroclor-1254	Sample et al. 1996
Arsenic	mouse	0.03	NOAEL	chronic	growth, reproduction	2.47			2.47	7.04E-01	2.34E+00	2.85E+00		EPA 2005 (Arsenic Eco SSL, geomean of TRVs for growth & reproduction)
	mouse	0.03	NOAEL	chronic	growth, reproduction	51.8			51.8	1.48E+01	4.90E+01	5.99E+01		EPA 2005 (Barium Eco SSL, geomean of TRVs
Barium														for growth & reproduction)
	mouse	0.03	NOAEL	chronic	growth, reproduction	18.0			18.0	_	_		HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
Benzo(b)fluoranthene										5.12E+00	1.70E+01	2.08E+01		growth & reproduction)
	mouse	0.03	NOAEL	chronic	growth, reproduction	18.0			18.0				HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
Benzo(g,h,i)perylene			110151							5.12E+00	1.70E+01	2.08E+01		growth & reproduction)
D (1)(1)	mouse	0.03	NOAEL	chronic	growth, reproduction	18.0			18.0	5 405 00	4 705 04	0.005.04	HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
Benzo(k)fluoranthene		0.00	NOAEL		111 4 114				7-	5.12E+00	1.70E+01	2.08E+01		growth & reproduction)
C11-C22 Aromatics	mouse	0.03	NOAEL	chronic	kidney toxicity	75			75	2.14E+01	7.10E+01	8.67E+01	Pyrene	MADEP 2003
C19-C36 Aliphatics	rat	0.35	NOEAL	chronic	liver toxicity	200			200	1.05E+02	3.50E+02	4.27E+02		MADEP 2003
C5-C8 Aliphatics	rat	0.35	LOAEL	chronic	neurotoxicity	407		10	40.7	2.14E+01	7.12E+01	8.70E+01	n-hexane	MADEP 2003
C9-C10 Aromatics	mouse	0.03	NOAEL	chronic	kidney toxicity	75			75	2.14E+01	7.10E+01	8.67E+01	Pyrene	MADEP 2003
	rat	0.35	NOAEL	chronic	liver weight	100			100				isoparaffins/napht	MADEP 2003
C9-C18 Aliphatics										5.27E+01	1.75E+02	2.14E+02	henes/n-alkanes	
	mouse	0.03	NOAEL	chronic	growth, reproduction	1.86			1.86	5.29E-01	1.76E+00	2.15E+00		EPA 2005 (Cadmium Eco SSL, geomean of TRVs
Cadmium						1.00				_				for growth & reproduction)
	mouse	0.03	NOAEL	chronic	growth, reproduction	2.40			2.40	6.85E-01	2.28E+00	2.78E+00		EPA 2005 (Chromium Eco SSL, geomean of
Chromium														TRVs for growth & reproduction)
	mouse	0.03	NOAEL	chronic	growth, reproduction	25.0			25.0	7.12E+00	2.36E+01	2.88E+01		EPA 2007 (Copper Eco SSL, geomean of TRVs
Copper			110151											for growth & reproduction)
5" (1) "	mouse	0.03	NOAEL	chronic	growth, reproduction	18.0			18.0	5 405 00	4 705 04	0.005.04	HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
Dibenz(a,h)anthracene		0.00	NOAFI	ala na mi a	anaceth named cation	40.7			40.7	5.12E+00	1.70E+01	2.08E+01		growth & reproduction)
Land	mouse	0.03	NOAEL	chronic	growth, reproduction	40.7			40.7	1.16E+01	3.86E+01	4.71E+01		EPA 2005 (Lead Eco SSL, geomean of TRVs for growth & reproduction)
Lead Mercury	mink	1	NOAEL	chronic	reproduction	1.0			1.0	6.85E-01	2.28E+00	2.78E+00		Sample et al. 1996
Mercury	mouse	0.03	NOAEL	chronic	growth, reproduction	7.70			7.70	2.19E+00	7.29E+00	8.90E+00		EPA 2007 (Nickel Eco SSL, geomean of TRVs for
Nickel	mouse	0.03	NOAEL	CHIOHIC	growin, reproduction	7.70			7.70	2.196+00	7.296+00	0.900+00		growth & reproduction)
Nickei	mouse	0.03	NOAEL	chronic	growth, reproduction	18.0			18.0				HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
Pyrene	mouse	0.03	NOAEL	CHIOTIC	growin, reproduction	10.0			10.0	5.12E+00	1.70E+01	2.08E+01	HIVIVV FAHS	growth & reproduction)
1 yiono	mouse	0.03	NOAEL	chronic	growth, reproduction	5.92			5.92	J. 12LT00	1.702701	2.00LT01	1	EPA 2005 (Vanadium Eco SSL, geomean of
Vanadium	mouse	0.00	I TOME	011101110	growin, reproduction	0.02			0.02	1.69E+00	5.61E+00	6.85E+00		TRVs for growth & reproduction)
- Graduiti	mouse	0.03	NOAEL	chronic	growth, reproduction	75.4			75.4	1.002100	0.012100	5.55E100		EPA 2007 (Zinc Eco SSL, geomean of TRVs for
Zinc	1110000	0.00	110/122	000	g. 5.7411, 10p1000001011	70.1			70.1	2.15E+01	7.14E+01	8.71E+01		growth & reproduction)
2.3.7.8-TCDD TEQ (mammalian)	rat	0.35	NOAEL	chronic	reproduction	0.000001	1		0.000001	5.27E-07	1.75E-06	2.14E-06	+	Sample et al. 1996

Table 4-5 Mammalian Toxicity Reference Values Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

LOAEL-Based TRVs

	Test Species	Test Species					Subchronic		Endpoint	Carnivorous	Herbivorous	Insectivorous		
		-								mammal	Mammal	Mammal		
	Common	Body	Endpoint	Study	Effect to	Toxicity	to Chronic		Adjusted	Red Fox	Deer Mouse	Short-tailed Shrew	Toxicity Value	Initial Compilation
Chemical	Name	Weight		Type	Test Organism	Value	UF		TRV	(mg/kg-day)	(mg/kg-day)	(mg/kg-day)	Surrogate	Source
		(kg)				(mg/kg-day)			(mg/kg-day)					
Antimony	mouse	0.03	LOAEL	chronic	growth, reproduction	2.76			2.76	7.86E-01	2.61E+00	3.19E+00		EPA 2005 (Antimony Eco SSL, geomean of TRVs
														for growth & reproduction)
Aroclor-1260	mouse	0.014	LOAEL	chronic	reproduction	0.68			0.68	1.60E-01	5.32E-01	6.50E-01	Aroclor-1254	Sample et al. 1996
	mouse	0.03	LOAEL	chronic	growth, reproduction	4.55			4.55	1.30E+00	4.31E+00	5.26E+00		EPA 2005 (Arsenic Eco SSL, geomean of TRVs
Arsenic					<u> </u>									for growth & reproduction)
Barium	mouse	0.03	LOAEL	chronic	growth, reproduction	82.7			82.7	2.36E+01	7.83E+01	9.55E+01		EPA 2005 (Barium Eco SSL, geomean of TRVs
														for growth & reproduction)
Benzo(b)fluoranthene	mouse	0.03	LOAEL	chronic	growth, reproduction	38.4			38.4				HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
										1.09E+01	3.64E+01	4.44E+01		growth & reproduction)
Benzo(g,h,i)perylene	mouse	0.03	LOAEL	chronic	growth, reproduction	38.4			38.4			=	HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
D (1)(1)		2.22	1015			00.4			20.4	1.09E+01	3.64E+01	4.44E+01	118484/ 50411	growth & reproduction)
Benzo(k)fluoranthene	mouse	0.03	LOAEL	chronic	growth, reproduction	38.4			38.4	4.005.04	0.045.04	4.445.04	HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
044 000 4		2.22	NOAEL		111 / 12	7.5		4.0	750	1.09E+01	3.64E+01	4.44E+01		growth & reproduction)
C11-C22 Aromatics	mouse	0.03	NOAEL	chronic	kidney toxicity	75		10	750	2.14E+02	7.10E+02	8.67E+02	Pyrene	MADEP 2003
C19-C36 Aliphatics	rat	0.35	NOAEL	chronic	liver toxicity	200		10	2000	1.05E+03	3.50E+03	4.27E+03	white mineral oils	MADEP 2003 MADEP 2003
C5-C8 Aliphatics	rat	0.35	LOAEL	chronic	neurotoxicity	407		4.0	407	2.14E+02	7.12E+02	8.70E+02	n-hexane	
C9-C10 Aromatics	mouse	0.03	NOAEL	chronic	kidney toxicity	75		10	750	2.14E+02	7.10E+02	8.67E+02		MADEP 2003
C9-C18 Aliphatics	rat	0.35	NOAEL	chronic	liver weight	100		10	1000	5.27E+02	4.755.00	2.14E+03	isoparaffins/napht	MADEP 2003
On desiran		0.00	1015	-11-		0.00			6.90	5.27E+02	1.75E+03	2.14E+03	henes/n-alkanes	EDA 0005 (O-designs E 00)
Cadmium	mouse	0.03	LOAEL	chronic	growth, reproduction	6.90			6.90	1.97E+00	6.54E+00	7.98E+00		EPA 2005 (Cadmium Eco SSL, geomean of TRVs
Ch va maio uma		0.03	LOAEL	ala u a usi a		58.2	+		58.2	1.97E+00	6.54E+00	7.98E+00		for growth & reproduction) EPA 2005 (Chromium Eco SSL, geomean of
Chromium	mouse	0.03	LOAEL	chronic	growth, reproduction	58.2			58.2	1.66E+01	5.51E+01	6.72E+01		TRVs for growth & reproduction)
Conner	maulaa	0.03	LOAFI		grouth reproduction					1.00=+01	3.31E+UI	0.725+01		
Copper	mouse	0.03	LOAEL	chronic	growth, reproduction	82.7			82.7	2.36E+01	7.83E+01	9.56E+01		EPA 2007 (Copper Eco SSL, geomean of TRVs for growth & reproduction)
Dibenz(a,h)anthracene	maulaa	0.03	LOAEL	chronic	growth, reproduction	38.4			38.4	2.30E+01	7.03E+01	9.300+01		EPA 2007 (PAH Eco SSL, geomean of TRVs for
Diberiz(a,ri)aritrilacerie	mouse	0.03	LOAEL	CHIOTIC	growin, reproduction	30.4			30.4	1.09E+01	3.64E+01	4.44E+01	HIVIVY PAHS	arowth & reproduction)
Lead	mouse	0.03	LOAEL	chronic	arowth, reproduction	186			186	1.032+01	3.04L+01	4.446101		EPA 2005 (Lead Eco SSL, geomean of TRVs for
Lead	IIIOuse	0.03	LOALL	CHIOTIC	growth, reproduction	100			100	5.31E+01	1.77E+02	2.15E+02		growth & reproduction)
Mercury	mink	1	NOAEL	chronic	reproduction	1.0		10	10.0	6.85E+00	2.28E+01	2.78E+01		Sample et al. 1996
Nickel	mouse	0.03	LOAEL	chronic	growth, reproduction	14.8		10	14.8	4.21E+00	1.40E+01	1.71E+01		EPA 2007 (Nickel Eco SSL, geomean of TRVs for
Nickei	mouse	0.03	LOALL	CHIOTIC	growth, reproduction	14.0			14.0	4.212+00	1.402+01	1.712+01		arowth & reproduction)
Pyrene	mouse	0.03	LOAEL	chronic	growth, reproduction	38.4			38.4	+			HMW PAHs	EPA 2007 (PAH Eco SSL, geomean of TRVs for
i yielle	mouse	0.03	LOALL	Cilionic	growth, reproduction	30.4			30.4	1.09E+01	3.64E+01	4.44E+01	TIWWTAIIS	growth & reproduction)
Zinc	mouse	0.03	LOAEL	chronic	growth, reproduction	298			298	8.48E+01	2.82E+02	3.44E+02		EPA 2007 (Zinc Eco SSL, geomean of TRVs for
	mouse	0.00	LONEL	011101110	growin, roproduction	200			200	0.402101	2.022102	0.446102		growth & reproduction)
2.3.7.8-TCDD TEQ (mammalian)	rat	0.35	LOAEL	chronic	reproduction	0.00001	+		0.00001	5.27E-06	1.75E-05	2.14E-05		Sample et al. 1996
=,0,.,0 . ODD 12 (1 100	0.00		011101110	TOPTOGGOGOT	0.00001	+ +		0.00001	J 0.2.7 E 00	1.702 00	2.112 00	+	1000

Notes:

mg - milligram

kg - kilogram LOAEL - lowest-observable-adverse-effect-level

NOAEL - no-observable-adverse-effect-level

TEF - toxicity equivalency factor

TRV - toxicity reference value

UF - uncertainty factor

Y - uncertainty factor used to derive TRV was included by original source

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Table 4-6
Avian Toxicity Reference Values
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

NOAEL-Based TRVs

NOAEL-Baseu IRVS										
Chemical	Test Species Common Name	Endpoint	Study Type	Effect to Test Organism	Toxicity Value (mg/kg-day)	LD ₅₀ to NOAEL UF	Subchronic to Chronic UF	Avian NOAEL-Equiv TRV (mg/kg-day)		Initial Compilation Source
Antimony	Black duck	NOAEL	chronic		1			1	Chromium	Sample et al. 1996
Aroclor-1260	Ring-Necked Pheasant	NOAEL	chronic	reproduction	0.18			0.18	Aroclor-1254	Sample et al. 1996
Arsenic	Chicken/Mallard duck	NOAEL	chronic	reproduction, growth, mortality	3.70			3.70049858		EPA 2005 (Arsenic Eco SSL, geomean of TRVs for growth, reproduction, & mortality)
Barium	1-day Old Chicks	NOAEL	sub-chronic	mortality	20.8			20.8		Sample et al. 1996
Benzo(a)anthracene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Benzo(a)Pyrene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Benzo(b)Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Benzo(g,h,i)perylene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Benzo(k)Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
C11-C22 Aromatics	Mallard duck	NOAEL	chronic		10000			10000	aromatic mixture	Patton and Dieter, 1979
C19-C36 Aliphatics	Mallard duck	NOAEL	chronic		400			400	paraffin	Patton and Dieter, 1979
C5-C8 Aliphatics	Mallard duck	NOAEL	chronic		400			400	paraffin	Patton and Dieter, 1979
C9-C10 Aromatics	Mallard duck	NOAEL	chronic		10000			10000	aromatic mixture	Patton and Dieter, 1979
C9-C18 Aliphatics	Mallard duck	NOAEL	chronic		400			400	paraffin	Patton and Dieter, 1979
Cadmium	various	NOAEL	chronic	growth, reproduction	1.47			1.467034136		EPA 2005 (Cadmium Eco SSL, geomean of TRVs for growth & reproduction)
Chromium	Chicken/Black duck/Turkey	NOAEL	chronic	growth, reproduction	2.66			2.657649427		EPA 2005 (Chromium Eco SSL, geomean of TRVs for growth & reproduction)
Chrysene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Copper	various	NOAEL	chronic	growth, reproduction	18.5			18.49433983		EPA 2007 (Copper Eco SSL, geomean of TRVs for growth & reproduction)
Dibenz(a,h)anthracene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	100		1.01	fluorene	Schafer et al. 1983
Lead	various	NOAEL	chronic	growth, reproduction	10.9			10.9408261		EPA 2005 (Lead Eco SSL, geomean of TRVs for growth & reproduction)
Mercury	Japanese Quail	NOAEL	chronic	reproduction	0.45			0.45		Sample et al. 1996
Nickel	Chicken/Duck	NOAEL	chronic	growth, reproduction	6.71			6.706541477		EPA 2007 (Nickel Eco SSL, geomean of TRVs for growth & reproduction)
Pyrene	Red-Winged Blackbird	LD50	acute	mortality	111	100		1.11	anthracene	Schafer et al. 1983
Vanadium	Chicken/Japanese quail	NOAEL	chronic	growth, reproduction	1.19			1.185971649		EPA 2005 (Vanadium Eco SSL, geomean of TRVs for growth & reproduction)
Zinc	various	NOAEL	chronic	growth, reproduction	66.1			66.06591637		EPA 2007 (Zinc Eco SSL, geomean of TRVs for growth & reproduction)
2,3,7,8-TCDD TEQ (avian)	Ring-Necked Pheasant	NOAEL	chronic	reproduction	1.40E-05		1	0.000014		Sample et al. 1996

Table 4-6
Avian Toxicity Reference Values
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

LOAEL-Based TRVs

LUAEL-Baseu IRVS										
Chemical	Test Species Common Name	Endpoint	Study Type	Effect to Test Organism	Toxicity Value (mg/kg-day)	LD ₅₀ to LOAEL UF	NOEAL to LOAEL UF	Avian NOAEL-Equiv TRV (mg/kg-day)	Toxicity Value Surrogate	Initial Compilation Source
Antimony	Black duck	LOAEL	chronic	reproduction	5			5	Chromium	Sample et al. 1996
Aroclor-1260	Ring-Necked Pheasant	LOAEL	chronic	reproduction	1.8			1.8	Aroclor-1254	Sample et al. 1996
Arsenic	Chicken/Mallard duck	LOAEL	chronic	reproduction, growth, mortality	4.51			4.51		EPA 2005 (Arsenic Eco SSL, geomean of TRVs for growth, reproduction, & mortality)
Barium	1-day Old Chicks	LOAEL	sub-chronic	mortality	41.7			41.7		Sample et al. 1996
Benzo(a)anthracene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Benzo(a)Pyrene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Benzo(b)Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Benzo(g,h,i)perylene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Benzo(k)Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
C11-C22 Aromatics	Mallard duck	NOAEL	chronic		10000		10	100000	aromatic mixture	Patton and Dieter, 1979
C19-C36 Aliphatics	Mallard duck	NOAEL	chronic		400		10	4000	paraffin	Patton and Dieter, 1979
C5-C8 Aliphatics	Mallard duck	NOAEL	chronic		400		10	4000	paraffin	Patton and Dieter, 1979
C9-C10 Aromatics	Mallard duck	NOAEL	chronic		10000		10	100000	aromatic mixture	Patton and Dieter, 1979
C9-C18 Aliphatics	Mallard duck	NOAEL	chronic		400		10	4000	paraffin	Patton and Dieter, 1979
Cadmium	various	LOAEL	chronic	growth, reproduction	6.35			6.35		EPA 2005 (Cadmium Eco SSL, geomean of TRVs for growth & reproduction)
Chromium	Chicken/Black duck	LOAEL	chronic	growth, reproduction	15.6			15.6		EPA 2005 (Chromium Eco SSL, geomean of TRVs for growth & reproduction)
Chrysene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Copper	various	LOAEL	chronic	growth, reproduction	34.9			34.9		EPA 2007 (Copper Eco SSL, geomean of TRVs for growth & reproduction)
Dibenz(a,h)anthracene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Fluoranthene	Red-Winged Blackbird	LD ₅₀	acute	mortality	101	10		10.1	fluorene	Schafer et al. 1983
Lead	various	LOAEL	chronic	growth, reproduction	44.6			44.6		EPA 2005 (Lead Eco SSL, geomean of TRVs for growth & reproduction)
Mercury	Japanese Quail	LOAEL	chronic	reproduction	0.9			0.9		Sample et al. 1996
Nickel	Chicken/Duck	LOAEL	chronic	growth, reproduction	18.6			18.6		EPA 2007 (Nickel Eco SSL, geomean of TRVs for growth & reproduction)
Pyrene	Red-Winged Blackbird	LD ₅₀	acute	mortality	111	10		11.1	anthracene	Schafer et al. 1983
Vanadium	Chicken/Japanese quail	LOAEL	chronic	growth, reproduction	1.70			1.70		EPA 2005 (Vanadium Eco SSL, geomean of TRVs for growth & reproduction)
Zinc	various	LOAEL	chronic	growth, reproduction	171			171		EPA 2007 (Zinc Eco SSL, geomean of TRVs for growth & reproduction)
2,3,7,8-TCDD TEQ (avian)	Ring-Necked Pheasant	LOAEL	chronic	reproduction	0.00014			0.00014		Sample et al. 1996

Notes:

mg - milligram kg - kilogram

LD₅₀ - lethal dose for 50% of study population LOAEL - lowest-observable-adverse-effect-level

NA - Not available

NOAEL - no-observable-adverse-effect-level

TEF - toxicity equivalency factor TRV - toxicity reference value

UF - uncertainty factor

Y - uncertainty factor used to derive TRV was included by original source

References

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Human Health and Enironmental Risk Characterization

		Location:	DD-N	IW-001	DD-N	IW-002	DD-MW-003	DD-MW-007							DD-MW-201				
		Sample ID:	DD-MW-001-R02-X	MW-1-051705	DD-MW-002-R01-X	DD-MW-002-R02-X	MW-03-051705	GZA-7-051705	DE	D-MW-201-001-D	DD-MW-201-001-X		DD-MW-201-R02-	Х	DD-MW-201-R04-	Х	DD-MW-201-R05-	Х	DD-MW-201-R06-D
		Sample Date:	6/5/2007	5/17/2005	6/5/2007	5/19/2008	5/17/2005	5/17/2005		6/26/2006	6/26/2006		8/14/2006		6/5/2007		12/11/2007		5/19/2008
		Depth:	11.6 - 16.6 feet	14 - 14 feet	9.8 - 14.8 feet	9.8 - 14.8 feet	3.5 - 3.5 feet	0 - 0 feet		9 - 9 feet	9 - 9 feet		10 - 10 feet		4 - 14 feet		4 - 14 feet		4 - 14 feet
CAS Number	Analyte	Únits	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q		Result Q	Result Q	2	Result	Q	Result	Q	Result	Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 UJ5	' 10 u'	*	5.1	UJ5	5.3	U	5.1	UJ	5.1 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA		0.32 u	0.33 u	1	0.3	u	0.32	U	0.3	U	0.31 U
50-32-8	BENZO(A)PYRENE	ug/l	0.2 U	NA	0.2 U	0.2 U	NA	NA		0.22 u	0.22 u	1	0.2	u	0.21	U	0.2	U	0.2 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA		0.32 U	0.33 U	J	0.3	U	0.32	U	0.3	U	0.31 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA		0.54 U	0.55 U	J	0.51	U	0.42	U	0.51	U	0.51 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA		0.32 U	0.33 U	J	0.3	U	0.32	U	0.3	U	0.31 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	5.1 J	NA	5.1 J	5.1 U	NA	NA		10 U	10 U	J	5.1	U	5.3	J	5.1	J	5.1 U
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 U	10 U	J	5.1	U	5.3	U	5.1	U	5.1 U
EPH1122	C11-C22 AROMATICS, ADJUSTED	D ug/l	100 U	NA	110 U	100 U	NA	NA		110 U	110 U	J	100	U	110	U	100	U	100 U
EPH1936	C19-C36 ALIPHATICS	ug/l	100 U	NA	110 U	100 U	NA	NA		110 U	110 U	J	100	U	110	U	100	U	100 U
PH912	C9-C18 ALIPHATICS	ug/l	100 U	NA	110 U	100 U	NA	NA		110 U	110 U	J	100	U	110	U	100	U	100 U
218-01-9	CHRYSENE	ug/l	1 U	NA	1 U	1 U	NA	NA		1.1 U	1.1 U	J	1	U	1.1	U	1	U	1 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA		0.54 U	0.55 U	J	0.51	U	0.42	U	0.51	U	0.51 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	5.1 J	NA	5.1 J	5.1 U	NA	NA		10 U	10 U	J	5.1	U	5.3	J	5.1	J	5.1 U
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 U	10 U	J	5.1	U	5.3	U	5.1	U	5.1 U
206-44-0	FLUORANTHENE	ug/l	1 U	NA	1 U	1 U	NA	NA		1.1 U	1.1 U	J	1	U	1.1	U	1	U	1 U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA		0.54 U	0.55 U	J	0.51	U	0.42	U	0.51	U	0.51 U
85-01-8	PHENANTHRENE	ug/l	0.2 U	NA	0.2 U	0.2 U	NA	NA		0.22	0.22 U	J	0.2	U	0.21	U	0.2	U	0.2 U
129-00-0	PYRENE	ug/l	1 U	NA	1.1 U	5.1 U	NA	NA		1.1 U	1.1 U	J	1	U	1.1	U	5.1	U	5.1 U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	13		10 U	10 U	J	10	U	1	U	1	U	1 U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	2 U	1 U	2 U	1 U	1 U	1 U		10 U	10 U	J	10	U	0.81	J	1	U	1 U
7440-39-3	BARIUM (DISSOLVED)	ug/l	47	39	32	19	34	110		40	40		42		44		38		45
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	1 U		1 U	1 U	J	1	U	1	U	2	U	1 U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	5.7		1 U	1 U	J	1	U	1	U	1	U	1 U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	35	1 U	7.3	1.9	43	1 U		5 U	5 U	J	5	U	25		22		3.5 J
7440-50-8	COPPER (DISSOLVED)	ug/l	NA	5 U	NA	NA	5 U	22		NA	NA		NA		NA		NA		NA
7439-92-1	LEAD (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	1 U		5 U	5 U	J	5	U	1	U	1	U	1 U
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U		0.2 U	0.2 U	J	0.2	U	0.2	J	0.71		0.2 U
7440-02-0	NICKEL (DISSOLVED)	ug/l	7.2	5.6	2.5	1.2	110	25		1.4 J	1.6 J	j	2.3	J	1.3		3.3		1.2
7782-49-2	SELENIUM (DISSOLVED)	ug/l	2 U	2 U	2 U	1 U	2 U	2 U		7.7 J	5.5 J	J	5.8	J	9.3	J	1	U	7.8
7440-22-4	SILER (DISSOLVED)	ug/l	1 U	0.1 U	1 U	1 U	0.1 U	0.1 U		5 U	5 U	J	1.1	J	1	U	1	U	1 U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 Ü	1 U	120		10 U	10 U	J	10	U	1	U	2	U	0.18 J
7440-66-6	ZINC (DISSOLVED)	ug/l	5 U	10 U	5 U	1.2 J	10 U	850		3.2 J	50 U	J	9.6	J	2.5	U	2.4	J	2.3 J
75-25-2	BROMOFORM	ug/l	NA	NA	NA NA	NA	NA	NA	i i	0.51 J	1 U	J	1	Ü	NA		NA		NA
108-88-3	TOLUENE	ua/l	NA	NA	NA	NA	NA	NA	i i	1 U	1 U	J	1	U	NA		NA		NA

Human Health and Enironmental Risk Characterization

	L	ocation:								DD-MW-203									DD-MW-204			
	Sa	mple ID:	DD-MW-201-R06	5-X	DD-MW-203-001	1-X	DD-MW-203-R02-X	DD-MW-203-R	03-X	DD-MW-203-R03-X-	FF_E	DD-MW-203-R04	4-X	DD-MW-203-R05-X		DD-MW-203-R06-X		DD-MW-204-001-X	DD-MW-204-R02-X	DD-MW-204-R05-)	х	DD-MW-205-001-X
	Samı	ple Date:	5/19/2008		6/26/2006		8/2/2006	8/14/2006		8/14/2006		6/5/2007		12/11/2007		5/19/2008		6/26/2006	6/5/2007	5/19/2008		6/26/2006
		Depth:	4 - 14 feet		8.5 - 8.5 feet		8.5 - 8.5 feet	10.5 - 10.5 fe	et	10.5 - 10.5 fee	t	4.5 - 12.5 feet	:	4.5 - 12.5 feet		4.5 - 12.5 feet		10.5 - 10.5 feet	7.5 - 13.5 feet	7.5 - 13.5 feet		9 - 9 feet
CAS Number	Analyte	Únits	Result	Q	Result	Q	Result Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result Q	Result C	Result	Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	5.2	U	10	UJ	NA	5.8	UJ	NA		5.2	U	5.1 L	JJ	5.1	U	10 U	5.1 L	0.84	J	10 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.31	U	0.94		NA	0.35	u	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
50-32-8	BENZO(A)PYRENE	ug/l	0.21	U	0.79		NA	0.23	u	0.23	u	0.2	U	0.2	U	0.2	U	0.21 U	0.2 L	0.2	U	0.2 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.31	U	0.74		NA	0.35	U	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.52	U	0.79		NA	0.58	U	0.57	u	0.4	U	0.51 U	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.31	U	0.76		NA	0.35	U	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.48	J	10	UJ	NA	5.8	U	NA		5.2	J	5.1	J	0.43	J	10 U	5.1	0.45	J	10 U
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	U	5.1 l	U	5.1	U	10 U	5.1 L	5.1	U	10 U
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 I	U	100	U	110 U	100 L	100	U	100 U
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 I	U	100	U	110 U	100 L	100	U	100 U
PH912	C9-C18 ALIPHATICS	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 I	U	100	U	110 U	100 L	100	U	100 U
218-01-9	CHRYSENE	ug/l	1	U	1	U	NA	1.2	U	1.1	u	1	U	1 1	U	1	U	1.1 U	1 L	1	U	1 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.52	U	0.82		NA	0.58	U	0.57	u	0.4	U	0.51 l	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	J	5.1	J	5.1	U	10 U	5.1	5.1	U	10 U
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	U	5.1 l	U	5.1	U	10 U	5.1 L	5.1	U	10 U
206-44-0	FLUORANTHENE	ug/l	1	U	1	U	NA	1.2	U	1.1	u	1	U	1 1	U	1	U	1.1 U	1 L	1	U	1 U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.52	U	0.81		NA	0.58	U	0.57	u	0.4	U	0.51 l	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
85-01-8	PHENANTHRENE	ug/l	0.21	U	0.23		NA	0.23	U	0.23	u	0.2	U	0.2	U	0.2	U	0.21 U	0.2 L	0.2	U	0.2 U
129-00-0	PYRENE	ug/l	5.2	U	1	U	NA	1.2	U	1.1	u	1	U	5.1 l	U	5.1	U	1.1 U	1 L	5.1	U	1 U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	1	U	10	U	NA	10	U	NA		1	U	1 1	U	1	U	10 U	1 L	1	U	10 U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	1	U	10	U	NA	3.9	J	NA		1	U	1 1	U	1	U	10 U	1 L	2	U	10 U
7440-39-3	BARIUM (DISSOLVED)	ug/l	46		14		NA	13		NA		12		11		11		26	17	20		7.3 J
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	NA	0.096	J	NA		1	U	2 1	U	1	U	0.16 J	1 L	1	U	1 U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1	U	1	U	NA	1	U	NA		1	U	1 1	U	1	U	1 U	1 L	1	U	1 U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	2.1	J	5	U	NA	5	U	NA		16		15		0.63	J	5 U	11	1	U	5 U
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA	NA		NA		NA		NA		NA		NA	NA	NA		NA
7439-92-1	LEAD (DISSOLVED)	ug/l	1	U	5	U	NA	5	U	NA		1	U	1	U	1	U	5 U	1 L	1	U	5 U
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	U	NA	0.2	U	NA		0.2	U	0.2	U	0.2	U	0.2 U	0.2 L	0.2	U	0.2 U
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.3		1.6	J	NA	1.3	J	NA		1.1		2	U	1.2		1.9 J	1.1	1.2		1.3 J
7782-49-2	SELENIUM (DISSOLVED)	ug/l	6.8		10	U	NA	7.3	J	NA		1	U	1 1	U	1	U	10 U	1 L	2	U	10 U
7440-22-4	SILER (DISSOLVED)	ug/l	1	U	5	U	NA	1.4	J	NA		1	U	1 1	U	1	U	5 U	1 L	1	U	5 U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	1	U	10	U	NA	10	U	NA		1	U	2	U	0.25	J	10 U	1 L	1	Ū	10 U
7440-66-6	ZINC (DISSOLVED)	ug/l	2.7		50	U	NA	1.4	J	NA		2.4	J	1.9	J	5.3		1.6 J	2.1	5	U	50 U
75-25-2	BROMOFORM	ug/l	NA		1	U	NA	1	U	NA		NA		NA		NA		0.67 J	NA	NA		1 U
108-88-3	TOLUENE	ua/l	NA		1	U	NA	1	U	NA		NA		NA		NA		1 U	NA	NA		1 U

Human Health and Enironmental Risk Characterization

	L	ocation:	C	D-MW	V-205									DD-MW-206									
	Sa	mple ID:	DD-MW-205-R02	-X	DD-MW-205-R0	03-X	DD-MW-205-R05-X	DD-MW-206-00	1-X	DD-MW-206-R02	2-D	DD-MW-206-R02	2-X	DD-MW-206-R03->	(DD-MW-206-R03-X-I	D	DD-MW-206-R04-X		DD-MW-206-R05-X	DD-MW-207-001-X	(DD-MW-207-R02-X
	Samp	le Date:	8/14/2006		6/6/2007		5/19/2008	6/26/2006		8/2/2006		8/2/2006		8/14/2006		8/14/2006		6/5/2007		5/19/2008	6/27/2006		8/2/2006
		Depth:	10 - 10 feet		5 - 11 feet		5 - 11 feet	8 - 8 feet		10 - 10 feet		10 - 10 feet		10 - 10 feet		10 - 10 feet		5 - 11 feet		5 - 11 feet	10 - 10 feet		12.5 - 12.5 feet
CAS Number	Analyte	Units	Result	Q	Result	Q	Result Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result Q	2	Result Q	Result	Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	UJ	5.1 l	UJ	5.2 U	J	5.1 U	10	U	NA
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
50-32-8	BENZO(A)PYRENE	ug/l	0.2	U	0.2	U	0.2 U	0.21	U	NA		NA		0.2	U	0.2	U	0.21 U	J	0.2 U	0.2	U	NA
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	10	U	5.1	J	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 J	J	0.49 J	10	U	NA
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 U	J	5.1 U	10	U	NA
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
PH912	C9-C18 ALIPHATICS	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
218-01-9	CHRYSENE	ug/l	1	U	1	U	1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	10	U	5.1	J	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 J	J	5.1 U	10	U	NA
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 U	J	5.1 U	10	U	NA
206-44-0	FLUORANTHENE	ug/l	1	U	1	U	1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
85-01-8	PHENANTHRENE	ug/l	0.2	U	0.2	U	0.2 U	0.21	U	NA		NA		0.2	U	0.2	U	0.21 U	J	0.2 U	0.2	U	NA
129-00-0	PYRENE	ug/l	1	U	1	U	5.1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	5.1 U	1	U	NA
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	10	U	1	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	10	U	NA
7440-38-2	ARSENIC (DISSOLVED)	ug/l	10	U	2	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	4.7	J	NA
7440-39-3	BARIUM (DISSOLVED)	ug/l	6.1	J	18		7.6	23		NA		NA		27		28		26		16	38		NA
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	1 U	1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1	U	1	U	1 U	1	U	NA		NA		1	U	1	U	1 U	J	1 U	0.69	J	NA
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	5	U	8.4		0.52 J	5	U	NA		NA		5	U	5	U	23		0.51 J	5	U	NA
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA	NA		NA		NA		NA		NA		NA		NA	NA		NA
7439-92-1	LEAD (DISSOLVED)	ug/l	5	U	1	U	1 U	5	U	NA		NA		5	U	5	U	1 U	J	1 U	5	U	NA
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	U	0.2 U	0.2	U	NA		NA		0.2	U	0.2	U	0.2 U	J	0.2 U	0.2	U	NA
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.3	J	1.4		0.59 J	1.4	J	NA		NA		1.3	J	10	U	0.91 J	J	0.69 J	4.7	J	NA
7782-49-2	SELENIUM (DISSOLVED)	ug/l	10	U	2	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	10	U	NA
7440-22-4	SILER (DISSOLVED)	ug/l	5	U	1	U	1 U	5	U	NA		NA		5	U	1.5	J	1 U	J	1 U	5	U	NA
7440-62-2	ANADIUM (DISSOLVED)	ug/l	10	U	1	U	0.29 J	10	Ū	NA		NA		10	U	10	Ü	1 U	J	1 U	10	Ū	NA
7440-66-6	ZINC (DISSOLVED)	ug/l	50	U	5	U	1.6 J	4.2	J	NA		NA		1.8	J	9.3	J	2.7		1.9 J	3.5	J	NA
75-25-2	BROMOFORM	ug/l	1	U	NA		NA	0.83	J	NA		NA		1	U	1	U	NA		NA	1	Ü	NA
108-88-3	TOLUENE	ua/l	1	U	NA		NA	1	U	NA		NA		1	U	0.52	J	NA		NA	1	U	NA

Human Health and Enironmental Risk Characterization

		Location:			DD-MW-207												DD-MW-208-R01-0	001-D				
		Sample ID:	DD-MW-207-R	03-X	DD-MW-207-R04	-X	DD-MW-207-R05	-D	DD-MW-207-R05	5-X	DD-MW-207-R0	6-X	DD-MW-208-R01-0	01-D	DD-MW-208-R01-0	01-X	DD-MW-208-R0	2-X	DD-MW-208-R0)3-X	DD-MW-208-R04	4-X
	Sa	ample Date:	8/14/2006		6/6/2007		12/11/2007		12/11/2007		5/19/2008		6/25/2007		6/25/2007		7/23/2007		12/11/2007		5/19/2008	
		Depth:	12.5 - 12.5 fe	eet	5 - 15 feet		5 - 15 feet		5 - 15 feet		5 - 15 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet	T
CAS Number	Analyte	Units	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q
51-28-5	2,4-DINITROPHENOL	ug/l	10	U	5.1	U	5.5	U	5.1	UJ	5.2	U	5.1	U	5.1	UJ	5.2	U	5.1	UJ	5.1	U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3	U	0.22	J	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
50-32-8	BENZO(A)PYRENE	ug/l	0.2	U	0.2	U	0.22	C	0.2	U	0.21	U	0.2	U	0.2	U	0.21	U	0.2	U	0.2	U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3	U	0.32	U	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.51	U	0.42	U	0.55	C	0.51	U	0.52	U	0.4	U	0.41	U	0.21	J	0.51	U	0.51	U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	10	U	5.1	J	5.5	J	5.1	J	0.51	J	5.1	U	5.1	U	0.78	J	5.1	J	1.6	J
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	10	U	5.1	U	5.5	U	5.1	U	5.2	U	5.1	U	5.1	U	0.34	J	5.1	U	5.1	U
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	110	U	100	C	100	U	100	U	290		460		100	U	NA		100	U
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	110	U	100	U	100	U	100	U	100	U	130		100	U	100	U	100	U
PH912	C9-C18 ALIPHATICS	ug/l	100	U	110	C	100	С	100	U	100	U	100	U	100	U	100	U	100	U	100	U
218-01-9	CHRYSENE	ug/l	1	U	0.24	J	1.1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.51	U	0.42	C	0.55	С	0.51	U	0.52	U	0.4	U	0.41	U	0.42	J	0.51	U	0.51	U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	10	U	5.1	J	5.5	J	5.1	J	5.2	U	5.1	J	5.1	J	5.2	J	5.1	J	1.5	J
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	10	U	0.39	J	5.5	U	5.1	U	5.2	U	5.1	U	5.1	U	5.2	U	5.1	U	5.1	U
206-44-0	FLUORANTHENE	ug/l	1	U	0.28	7	1.1	C	1	U	1	U	1.885	J	1.875	J	1	U	1	U	1	U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.51	U	0.42	U	0.55	U	0.51	U	0.52	U	0.4	U	0.41	U	0.38	J	0.51	U	0.51	U
85-01-8	PHENANTHRENE	ug/l	0.2	U	0.2	U	0.22	C	0.2	U	0.21	U	0.62	J	0.37	J	0.21	U	0.2	U	0.2	U
129-00-0	PYRENE	ug/l	1	U	0.29	J	5.5	U	5.1	U	5.2	U	0.28	J	1	U	1	U	5.1	U	5.1	U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	2.3	U	1	C	1	С	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	7.7	J	5.7		10		10		3.9		1.7	J	1.7	J	NA		1	U	1	U
7440-39-3	BARIUM (DISSOLVED)	ug/l	44		38		44		45		31		27		26		NA		25		31	
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	2	U	2	U	1	U	2	U	1	U	NA		2	U	1	U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	0.84	J	1	C	1	С	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	5	U	4.9		9.1		10		1.3		9.8		11		NA		25		0.36	J
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA		NA		NA		NA		NA		NA		NA		NA	
7439-92-1	LEAD (DISSOLVED)	ug/l	5	U	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	0.46	J
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	C	0.2	С	0.2	U	0.2	U	0.2	U	0.2	U	NA		0.2	U	0.2	U
7440-02-0	NICKEL (DISSOLVED)	ug/l	4	J	1	U	2	U	2	U	0.43	J	3.6		3.2		NA		0.96	J	2	
7782-49-2	SELENIUM (DISSOLVED)	ug/l	10	U	1	U	1	U	1	U	1	U	2	U	2	U	NA		1	U	1	U
7440-22-4	SILER (DISSOLVED)	ug/l	5	U	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	0.92	J	1	U	2	U	2	U	1	U	0.91	J	1	J	NA		2	U	0.52	J
7440-66-6	ZINC (DISSOLVED)	ug/l	3.5	J	2.5	U	1.2	J	1.5	J	1.9	J	9.4		7.7		NA		20		110	
75-25-2	BROMOFORM	ug/l	1	U	NA		NA		NA		NA		5	U	5	U	NA		NA		NA	
108-88-3	TOLUENE	ug/l	1	U	NA		NA		NA		NA		5	U	5	U	NA		NA		NA	T

Table 4-7
Estimated Potential Hazard Quotients for Ecological Receptors in DDA
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

		orous		vorous	Herbiy			orous	Insecti			orous
Analyte	Man			nmal	Man			ian		ian		ian
, 	Red	-		led Shrew	Meado			ed Hawk		lcock		ıail
	NOAEL HQ						NOAEL HQ				NOAEL HQ	
Antimony	3E-06	2E-05	4E-03	2E-02	1E-03	6E-03	6E-04	1E-04	1E-01	2E-02	5E-04	1E-04
Arsenic	4E-05	2E-05	9E-02	5E-02	1E-02	8E-03	4E-04	3E-04	9E-02	7E-02	3E-04	3E-04
Barium	8E-05	5E-05	1E-01	6E-02	6E-02	4E-02	3E-03	2E-03	5E-01	3E-01	6E-03	3E-03
Cadmium	4E-05	1E-05	7E-01	2E-01	2E-02	4E-03	4E-04	8E-05	8E-01	2E-01	8E-04	2E-04
Chromium	3E-03	1E-04	2E+00	6E-02	6E-01	3E-02	3E-02	4E-03	3E+00	5E-01	2E-02	3E-03
Copper	5E-05	2E-05	1E-01	3E-02	3E-02	9E-03	9E-04	5E-04	2E-01	1E-01	1E-03	8E-04
Lead	8E-05	2E-05	2E-01	3E-02	2E-02	4E-03	3E-03	8E-04	8E-01	2E-01	2E-03	6E-04
Mercury	3E-06	3E-07	3E-02	3E-03	4E-03	4E-04	3E-04	1E-04	2E-01	9E-02	8E-04	4E-04
Nickel	8E-04	4E-04	4E-01	2E-01	2E-01	1E-01	1E-02	4E-03	1E+00	5E-01	8E-03	3E-03
Vanadium	1E-04	7E-05	2E-01	1E-01	4E-02	3E-02	9E-03	6E-03	2E+00	1E+00	6E-03	4E-03
Zinc	4E-04	9E-05	4E-01	9E-02	6E-02	1E-02	2E-03	9E-04	4E-01	2E-01	3E-03	1E-03
Aroclor-1260	4E-05	4E-06	1E+00	1E-01	2E-02	2E-03	2E-04	2E-05	3E-01	3E-02	2E-04	2E-05
2,3,7,8-TCDD TEQ (mammalian)	2E-04	2E-05	9E+00	9E-01	9E-02	9E-03	NA	NA	NA	NA	NA	NA
2,3,7,8-TCDD TEQ (avian)	NA	NA	NA	NA	NA	NA	9E-04	9E-05	2E+00	2E-01	7E-04	7E-05
Benzo(b)fluoranthene	3E-07	1E-07	7E-03	3E-03	3E-04	1E-04	8E-05	8E-06	1E-01	1E-02	2E-04	2E-05
Benzo(g,h,i)perylene	2E-07	8E-08	5E-03	2E-03	2E-04	9E-05	5E-05	5E-06	9E-02	9E-03	1E-04	1E-05
Benzo(k)fluoranthene	3E-07	1E-07	7E-03	3E-03	2E-04	9E-05	8E-05	8E-06	1E-01	1E-02	1E-04	1E-05
Dibenz(a,h)anthracene	2E-07	8E-08	4E-03	2E-03	1E-04	6E-05	5E-05	5E-06	7E-02	7E-03	8E-05	8E-06
Pyrene	5E-07	3E-07	1E-02	5E-03	1E-03	6E-04	2E-04	2E-05	2E-01	2E-02	8E-04	8E-05
C5-C8 Aliphatics	6E-06	6E-07	4E-03	4E-04	1E-03	1E-04	6E-06	6E-07	9E-04	9E-05	1E-05	1E-06
C9-C10 Aromatics	2E-05	2E-06	1E-02	1E-03	3E-03	3E-04	6E-07	6E-08	9E-05	9E-06	1E-06	1E-07
C9-C18 Aliphatics	2E-06	2E-07	1E-03	1E-04	3E-04	3E-05	4E-06	4E-07	6E-04	6E-05	7E-06	7E-07
C11-C22 Aromatics	9E-05	9E-06	6E-02	6E-03	2E-02	2E-03	4E-06	4E-07	5E-04	5E-05	6E-06	6E-07
C19-C36 Aliphatics	3E-05	3E-06	2E-02	2E-03	6E-03	6E-04	2E-04	2E-05	2E-02	2E-03	2E-04	2E-05

Bold indicates HQ exceeds threshold of 1

HQ - Hazard quotient

NOAEL - No observed adverse effect level

LOAEL - Lowest observed adverse effect level

Attachment A (1) Analytical Data - DDA Soil Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

	1	Location	DD-GP-003	DD-GP-004	DD-GP-006	DD-GP-009	DD-GP-207	DD-SB-206	DD-SS-003	DD-SS-005	DD-SS-007	DD-SS-012	DD-SS-014	DD-T	P-013
			DDA-GP-3 RR & Dup	DD-GF-004 DDA-GP-4	DDA-GP-6	DDA-GP-9 RR & Dup		DD-SB-206-003-X	DD-SS-001-001-X	DD-SS-005	DD-SS-007-001-X & Dup	DD-SS-012-001-X	DD-SS-014-001-X	DD-TP-001-001-X	DD-TP-001-002-X
		Sample Date	5/16/2005	5/16/2005	5/16/2005	5/16/2005	6/9/2006	6/19/2006	12/8/2005	9/21/2006	10/26/2006	5/30/2007	5/30/2007	11/21/2005	11/21/2005
CAS		Depth	0 - 3 feet	2 - 4 feet	0 to 3 feet	0 to 3 feet	0 to 2 feet	0 to 2 feet	0 to 3 feet	0 to 0.5 feet	0 to 0.5 feet	0 to 2 feet	0 to 2 feet	0 to 2 feet	2 to 4 feet
Number	Analyte	Units	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q
7440-36-0	ANTIMONY	mg/kg	1.30 U	1.65 U	1.35 U	1.15 U	1.30 J	NA	NA	1.70	NA	NA	NA	NA	NA
7440-38-2	ARSENIC	mg/kg	9.5	2.40	3.40	2.60	4.10 J	NA	NA	4.20	NA	NA	NA	NA	NA
7440-39-3	BARIUM	mg/kg	52	43.00	220.00	24.00	59.00	NA	NA	47.00	NA	NA	NA	NA	NA
7440-41-7	BERYLLIUM	mg/kg	0.13 U	0.17 U	0.55	0.12 U	0.65 U	NA	NA	0.13 U	NA	NA	NA	NA	NA
7440-43-9	CADMIUM	mg/kg	0.61	0.43	0.78	0.38	1.40	NA	NA	1.30	NA	NA	NA	NA	NA
7440-47-3	CHROMIUM	mg/kg	28	11.00	230.00	16.00	17.00	NA NA	NA	34.00	NA NA	NA	NA	NA	NA
7440-50-8	COPPER	mg/kg	NA 110	NA 22.22	NA	NA 40.00	NA 450.00	NA	NA	56.00	NA NA	NA	NA	NA	NA
7439-92-1	LEAD	mg/kg	110	20.00	60.00	40.00	150.00	NA NA	NA	88.00	NA NA	NA NA	NA NA	NA NA	NA NA
7439-97-6 7440-02-0	MERCURY NICKEL	mg/kg mg/ka	NA 20	0.062 26.00	0.13 190.00	0.081 20.00	0.29 18.00	NA NA	NA NA	2.90 35.00	NA NA	NA NA	NA NA	NA NA	NA NA
7440-02-0	SILVER	3 3	1.4	0.85 U	0.65 U	0.55 U	18.00 1.60 U	NA NA	NA NA	0.83	NA NA	NA NA	NA NA	NA NA	NA NA
7440-22-4	VANADIUM	mg/kg mg/kg	23	22.00	38.00	9.30	26.00	NA NA	NA NA	23.00	NA NA	NA NA	NA NA	NA NA	NA NA
7440-62-2	ZINC	mg/kg	140	68.00	130.00	75.00	88.00	NA NA	NA NA	23.00 NA	NA NA	NA NA	NA NA	NA NA	NA NA
11096-82-5	AROCLOR-1260	mg/kg	0.055 U	0.065 U	0.055 U	0.055 U	0.055 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	0.060 U	0.060 U
95-63-6	1.2.4-TRIMETHYLBENZENE	mg/kg	0.033 0 NA	NA NA	NA NA	NA NA	0.0039 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA	0.000 C
107-06-2	1.2-DICHLOROETHANE	mg/kg	NA NA	NA NA	NA NA	NA NA	0.0019 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
108-67-8	1.3.5-TRIMETHYLBENZENE	mg/kg	NA NA	NA NA	NA NA	NA NA	0.0019 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
99-87-6	4-ISOPROPYLTOLUENE	mg/kg	NA NA	NA NA	NA	NA NA	0.040	NA NA	NA NA	NA NA	NA NA	NA NA	NA	NA NA	NA
67-64-1	ACETONE	mg/kg	NA NA	NA NA	NA	NA NA	0.24 J	NA NA	NA NA	NA NA	NA	NA NA	NA.	NA NA	NA NA
91-20-3	NAPHTHALENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.019 U	NA NA	0.11 U	0.27	NA NA	NA NA	NA NA	NA NA	NA NA
104-51-8	N-BUTYLBENZENE	ma/ka	NA NA	NA	NA	NA NA	NA	NA	NA	NA	NA NA	NA	NA	NA	NA
103-65-1	N-PROPYLBENZENE	mg/kg	NA	NA	NA	NA	0.0019 U	NA	NA	NA	NA	NA	NA	NA	NA
98-06-6	TERT-BUTYLBENZENE	mg/kg	NA	NA	NA	NA	0.0019 U	NA	NA	NA	NA	NA	NA	NA	NA
108-88-3	TOLUENE	mg/kg	NA	NA	NA	NA	0.14 J	NA	NA	NA	NA	NA	NA	NA	NA
91-57-6	2-METHYLNAPHTHALENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.066	NA	NA	NA	NA	NA
106-47-8	4-CHLOROANILINE	mg/kg	0.385 U	0.13 J	0.11 J	0.37 U	0.39 U	NA	0.41 U	NA	NA	NA	NA	NA	NA
56-55-3	BENZO(A)ANTHRACENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.13	NA	NA	NA	NA	NA
50-32-8	BENZO(A)PYRENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.095 U	NA	NA	NA	NA	NA
205-99-2	BENZO(B)FLUORANTHENE	mg/kg	0.095 U	0.11 U	0.090 U	0.14 J	0.10 U	NA	0.11 U	0.17	NA	NA	NA	NA	NA
191-24-2	BENZO(G,H,I)PERYLENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.095 U	NA	NA	NA	NA	NA
207-08-9	BENZO(K)FLUORANTHENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.11 U	NA	0.11 U	0.11 U	NA	NA	NA	NA	NA
218-01-9	CHRYSENE	mg/kg	0.095 U	0.11 U	0.090 U	0.14 J	0.10 U	NA	0.11 U	0.16	NA	NA	NA	NA	NA
53-70-3	DIBENZ(A,H)ANTHRACENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.095 U	NA	NA	NA	NA	NA
206-44-0	FLUORANTHENE	mg/kg	0.095 U	0.11 U	0.090 U	0.16 J	0.38	NA	0.11 U	0.19	NA	NA	NA	NA	NA
193-39-5	INDENO(1,2,3-CD)PYRENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.19 U	NA	0.11 U	0.20 U	NA	NA	NA	NA	NA
85-01-8	PHENANTHRENE	mg/kg	0.095 U	0.11 U	0.090 U	0.17 J	0.38	NA	0.11 U	0.17	NA	NA	NA	NA	NA
129-00-0	PYRENE	mg/kg	0.095 U	0.11 U	0.090 U	0.21	0.34	NA	0.11 U	0.22	NA	NA	NA	NA	NA
83-32-9	ACENAPHTHENE	mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA	0.11 U	0.095 U	NA NA	NA	NA	NA	NA
EPH1122	C11-C22 AROMATICS, ADJUSTED	mg/kg	55 170	17.00 15.00	14.00	54.00	15.00	NA NA	17.00 V 110.00 V	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
EPH1936	C19-C36 ALIPHATICS	mg/kg			23.00	120.00	5.30					NA NA		NA NA	NA NA
VPH912	C9-C18 ALIPHATICS	mg/kg	1.95 U	2.15 U	1.90 U	6.30	1.95 U	NA NA	19.00 V	NA NA	NA NA		NA NA		
VPH58 VPH910	C5-C8 ALIPHATICS, ADJUSTED C9-C10 AROMATICS	mg/kg	NA NA	NA NA	NA NA	NA NA	2.75 U 2.75 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
		mg/kg	NA NA	NA NA		NA NA	0.0019 U		NA NA			NA NA		NA NA	NA NA
100-41-4	ETHYLBENZENE O VVI ENE	mg/kg	NA NA	NA NA	NA NA	NA NA	0.0019 U	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
95-47-6	O-XYLENE ANTHRACENE	mg/kg	0.095 U	0.11 U	0.090 U	0.22	0.0019 U 0.10 U	NA NA	0.11 U	0.095 U	NA NA	NA NA	NA NA	NA NA	NA NA
120-12-7 86-73-7	FLUORENE	mg/kg mg/kg	0.095 U	0.11 U 0.11 U	0.090 U	0.22 0.095 U	0.10 U	NA NA	0.11 U	0.095 U	NA NA	NA NA	NA NA	NA NA	NA NA
00-13-1	M.P-XYLENES	mg/kg mg/kg	0.095 U	0.11 U	0.090 U	0.095 U	0.10 U	NA NA	0.11 U	0.095 0	NA NA	NA NA	NA NA	NA NA	NA NA
1746-01-6	2,3,7,8-TCDD TEQ	mg/kg mg/kg	1.50E-05	1.30E-06	4.80E-06	7.50E-07	1.10E-06	9.23E-05	NA NA	NA	2.80E-05	1.40E-06	6.20E-05	NA NA	NA NA
1740-01-6	[2,3,1,0-10DD 1EQ	mg/kg	1.5UE-U5	1.3UE-U6	4.6UE-U0	/ .5UE-U/	1.10E-06	9.23E-U5	INA	INA	∠.8UE-U5	1.40⊑-06	0.ZUE-U5	INA	INA

Notes:
DDA - Demolition Debris Area
mg/kg - milligram per kilogram
U - non-detect
J - concentration is an estimated value
NA - not analyzed

Q - qualifier
Only detected analytes are presented herein

1 of 2 11/2/2011 Attachment A (1) Analytical Data - DDA Soil Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

		Location	DD-TP-003	DD-TP-004	DD-TP-005	DD-TP-007	DD-TP-008	DD-TP-009	DD-TP-012	DD-TP-201	DD-TI	P-202
		Sample ID	DD-11-003 DD-TP03-2S	DD-TP-004 DD-TP04-2S	DD-TP-003	DD-17-007 DD-TP07-2S-D & Dup	DD-TP-000 DD-TP08-2S	DD-17-009 DD-TP09-2S-D & Dup	DD-TP12-5W	DD-TP-201-001-X	DD-TP-202-002-X	DD-TP-202-003-X
		Sample Date	12/20/2004	12/20/2004	12/20/2004	12/20/2004	12/20/2004	12/20/2004	12/20/2004	6/6/2006	6/28/2006	6/28/2006
CAS		Depth	0 to 3 feet	0 to 3 feet	0 to 3 feet	0 to 3 feet	0 to 3 feet	0 to 3 feet	0 to 3 feet	1 to 2 feet	2 to 3 feet	2 to 3 feet
Number	Analyte	Units	Result Q	Result Q	Result Q		Result Q		Result Q		Result Q	Result Q
7440-36-0	ANTIMONY	mg/kg	3.10	1.35 U	8.10	NA	1.20 U	NA	3.90 U	NA	1.50 J	1.50 U
7440-38-2	ARSENIC	mg/kg	26.00	0.65 U	7.60	4.00	1.30	1.92	1.95 U	2.10 J	2.80 J	2.95 U
7440-39-3	BARIUM	mg/kg	79.00	18.00	240.00	71.00	19.00	1395.00	1400.00	500.00	30.00	1.70 J
7440-41-7	BERYLLIUM	mg/kg	0.15 U	0.14 U	0.73	0.14 U	0.12 U	0.14 U	4.10	0.65 U	0.60 U	0.60 U
7440-43-9	CADMIUM	mg/kg	0.50	0.14 U	2.60	0.94	0.12 U	1.75	2.10	1.40	0.57 J	0.60 U
7440-47-3 7440-50-8	CHROMIUM COPPER	mg/kg	130.00 NA	8.30 NA	2200.00 NA	45.50 NA	9.80 NA	24.50 NA	300.00 NA	120.00 NA	29.00 NA	1.20 J NA
7439-92-1	LEAD	mg/kg mg/kg	300.00	9.50	450.00	145.00	11.00	57.00	73.00	77.00	110.00	3.00
7439-92-1	MERCURY	mg/kg	0.40	9.50 0.055 U	0.42	0.75	0.050 U	1.20	0.055 U	0.21	0.030 U	0.021 U
7440-02-0	NICKEL	mg/kg	1400.00	7.80	1200.00	NA NA	11.00	NA NA	260.00	95.00	74.00	0.021 J
7440-02-0	SILVER	mg/kg	0.75 U	0.65 U	3.20	0.70 U	0.60 U	0.65 U	1.95 U	1.70 U	1.50 U	1.50 U
7440-62-2	VANADIUM	mg/kg	29.00	17.00	150.00	35.50	13.00	27.50	120.00	71.00	16.00	1.30 J
7440-66-6	ZINC	mg/kg	150.00	34.00	810.00	195.00	45.00	175.00	510.00	130.00	110.00	4.10 J
11096-82-5	AROCLOR-1260	mg/kg	0.060 U	0.050 U	0.055 U	1.35	0.055 U	0.055 U	0.065 U	0.055 U	0.055 U	NA NA
95-63-6	1,2,4-TRIMETHYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	2.00	NA
107-06-2	1,2-DICHLOROETHANE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.0020	0.0016 U	0.0013 U	0.14 U	NA
108-67-8	1,3,5-TRIMETHYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	NA	0.0016 U	0.0013 U	0.68	NA
99-87-6	4-ISOPROPYLTOLUENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.14 U	NA
67-64-1	ACETONE	mg/kg	NA	0.12 U	NA	NA	0.11 U	0.095 U	0.16 U	0.13 U	13.50 U	NA
91-20-3	NAPHTHALENE	mg/kg	0.24 J	0.012 U	0.19 U	0.21 U	0.011 U	0.0095 U	0.016 U	0.29	0.095 U	NA
104-51-8	N-BUTYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.13 J	NA
103-65-1	N-PROPYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.28	NA
98-06-6	TERT-BUTYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.24 J	NA
108-88-3	TOLUENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.14 J	NA
91-57-6	2-METHYLNAPHTHALENE	mg/kg	0.20 U	0.18 U	0.19 U	0.21 U	0.18 U	0.19 U	0.11 U	0.19 U	0.22	NA
106-47-8	4-CHLOROANILINE	mg/kg	0.80 U	0.70 U	1.50 U	0.85 U	0.70 U	0.75 U	0.42 U	1.90 U	0.38 U	NA
56-55-3	BENZO(A)ANTHRACENE	mg/kg	0.35	0.52	0.34 J	0.21 U	0.18 U	0.19 U	0.11 U	1.59	0.095 U	NA
50-32-8	BENZO(A)PYRENE	mg/kg	0.31	0.79	0.41	0.21 U	0.18 U	0.19 U	0.11 U	1.44	0.095 U	NA
205-99-2	BENZO(B)FLUORANTHENE	mg/kg	0.20 U	0.52	0.19 U	0.21 U	0.18 U	0.19 U	0.11 U	2.30	0.095 U	NA
191-24-2	BENZO(G,H,I)PERYLENE	mg/kg	0.20 U	0.18 U	0.19 U	0.21 U	0.18 U	0.19 U	0.11 U	1.30	0.095 U	NA NA
207-08-9	BENZO(K)FLUORANTHENE	mg/kg	0.31	0.69	0.42	0.21 U	0.18 U	0.19 U	0.11 U	2.10	0.10 U	NA NA
218-01-9 53-70-3	CHRYSENE DIBENZ(A.H)ANTHRACENE	mg/kg ma/ka	0.44 J 0.20 U	0.94 0.18 U	0.60 J 0.19 U	0.22 U 0.21 U	0.18 U 0.18 U	0.19 U 0.19 U	0.11 U 0.11 U	1.69 1.30	0.095 U 0.095 U	NA NA
206-44-0	FLUORANTHENE	mg/kg	0.20 0	1.99	1.17	0.21 0	0.18 U	0.19 U	0.11 U	4.40	0.095 U	NA NA
193-39-5	INDENO(1.2.3-CD)PYRENE	mg/kg	0.95 0.20 U	0.18 U	0.19 U	0.41 0.21 U	0.18 U	0.19 U	0.11 U	0.94	0.095 U	NA NA
85-01-8	PHENANTHRENE	mg/kg	0.20 U	1.84	0.19 0	0.21 U	0.18 U	0.19 U	0.11 U	4.35	0.19 U	NA NA
129-00-0	PYRENE	mg/kg	1.075	1.89	1.050	0.25 U	0.18 U	0.19 U	0.11 U	4.050	0.095 U	NA NA
83-32-9	ACENAPHTHENE	mg/kg	0.20 U	0.18 U	0.19 U	0.34 J 0.21 U	0.18 U	0.19 U	0.11 U	0.46	0.095 U	NA NA
EPH1122	C11-C22 AROMATICS. ADJUSTED	mg/kg	29.00	35.00	22.00	NA NA	1.85 U	11.50	69.00	57.00	18.00	NA NA
EPH1936	C19-C36 ALIPHATICS	mg/kg	20.00	32.00	56.00	NA NA	1.85 U	20.00	55.00	19.00	72.00	NA NA
VPH912	C9-C18 ALIPHATICS	mg/kg	2.00 U	4.80	1.90 U	2.15 U	1.85 U	1.90 U	2.050 U	1.85 U	11.00	NA
VPH58	C5-C8 ALIPHATICS, ADJUSTED	mg/kg	NA NA	1.25 U	NA NA	NA NA	0.90 U	0.75 U	NA NA	1.40 U	5.20	NA
VPH910	C9-C10 AROMATICS	mg/kg	NA	1.25 U	NA	NA NA	0.90 U	0.75 U	NA	1.40 U	24.00	NA
100-41-4	ETHYLBENZENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.23	NA
95-47-6	O-XYLENE	mg/kg	NA	0.0012 U	NA	NA	0.0011 U	0.00095 U	0.0016 U	0.0013 U	0.49	NA
120-12-7	ANTHRACENE	mg/kg	0.20 U	0.64	0.19 U	0.21 U	0.18 U	0.19 U	0.11 U	1.035	0.095 U	NA
86-73-7	FLUORENE	mg/kg	0.20 U	0.18 U	0.19 U	0.21 U	0.18 U	0.19 U	0.11 U	0.74	0.095 U	NA
	M,P-XYLENES	mg/kg	0.27 J	0.18 U	0.38 U	0.21 U	0.18 U	0.19 U	0.11 U	0.47 U	0.095 U	NA
1746-01-6	2,3,7,8-TCDD TEQ	mg/kg	9.90E-06	NA	9.50E-05	1.90E-05	NA	NA	NA	NA	NA	NA

Notes:
DDA - Demolition Debris Area
mg/kg - milligram per kilogram
U - non-detect
J - concentration is an estimated value
NA - not analyzed

Q - qualifier
Only detected analytes are presented herein

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Human Health and Enironmental Risk Characterization

		Location:	DD-N	IW-001	DD-N	IW-002	DD-MW-003	DD-MW-007						DD-MW-201			
		Sample ID:	DD-MW-001-R02-X	MW-1-051705	DD-MW-002-R01-X	DD-MW-002-R02-X	MW-03-051705	GZA-7-051705	DD-M	W-201-001-D	DD-MW-201-001-X	DD-MW-2	201-R02-X	DD-MW-201-R04	I-X DD-N	IW-201-R05-X	DD-MW-201-R06-D
		Sample Date:	6/5/2007	5/17/2005	6/5/2007	5/19/2008	5/17/2005	5/17/2005	6/	/26/2006	6/26/2006	8/14/	/2006	6/5/2007	1	2/11/2007	5/19/2008
		Depth:	11.6 - 16.6 feet	14 - 14 feet	9.8 - 14.8 feet	9.8 - 14.8 feet	3.5 - 3.5 feet	0 - 0 feet	9) - 9 feet	9 - 9 feet	10 - 1	0 feet	4 - 14 feet	4	1 - 14 feet	4 - 14 feet
CAS Number	Analyte	Units	Result Q	Result Q	Result Q	Result Q	Result Q	Result Q	Re	esult Q	Result C	Resu	lt Q	Result	Q R	esult Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 UJ5	5* 10 u	5.1	UJ:	5 5.3	U	5.1 U.	5.1 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA	0).32 u	0.33 ι	0.3	u	0.32	U	0.3 U	0.31 U
50-32-8	BENZO(A)PYRENE	ug/l	0.2 U	NA	0.2 U	0.2 U	NA	NA	0).22 u	0.22 u	0.2	u	0.21	U	0.2 U	0.2 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA	0).32 U	0.33 L	0.3	U	0.32	U	0.3 U	0.31 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA	0).54 U	0.55 L	0.51	U	0.42	U	0.51 U	0.51 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3 U	NA	0.3 U	0.3 U	NA	NA	0).32 U	0.33 L	0.3	U	0.32	U	0.3 U	0.31 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	= ug/l	5.1 J	NA	5.1 J	5.1 U	NA	NA		10 U	10 L	5.1	U	5.3	J	5.1 J	5.1 U
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 U	10 L	5.1	U	5.3	U	5.1 U	5.1 U
EPH1122	C11-C22 AROMATICS, ADJUSTED	D ug/l	100 U	NA	110 U	100 U	NA	NA	1	110 U	110 L	100	U	110	U	100 U	100 U
EPH1936	C19-C36 ALIPHATICS	ua/l	100 U	NA	110 U	100 U	NA	NA	1	110 U	110 L	100	U	110	U	100 U	100 U
PH912	C9-C18 ALIPHATICS	ug/l	100 U	NA	110 U	100 U	NA	NA	1	110 U	110 L	100	U	110	U	100 U	100 U
218-01-9	CHRYSENE	ua/l	1 U	NA	1 U	1 U	NA	NA		1.1 U	1.1 L	1	U	1.1	U	1 U	1 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA	0).54 U	0.55 L	0.51	U	0.42	U	0.51 U	0.51 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	5.1 J	NA	5.1 J	5.1 U	NA	NA	,	10 U	10 L	5.1	U	5.3	J	5.1 J	5.1 U
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	5.1 U	NA	5.1 U	5.1 U	NA	NA		10 U	10 L	5.1	U	5.3	U	5.1 U	5.1 U
206-44-0	FLUORANTHENE	ug/l	1 U	NA	1 U	1 U	NA	NA		1.1 U	1.1 L	1	U	1.1	U	1 U	1 U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.4 U	NA	0.42 U	0.51 U	NA	NA	0).54 U	0.55 L	0.51	U	0.42	U	0.51 U	0.51 U
85-01-8	PHENANTHRENE	ug/l	0.2 U	NA	0.2 U	0.2 U	NA	NA	0).22	0.22 L	0.2	U	0.21	U	0.2 U	0.2 U
129-00-0	PYRENE	ug/l	1 U	NA	1.1 U	5.1 U	NA	NA		1.1 U	1.1 L	1	U	1.1	U	5.1 U	5.1 U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	13		10 U	10 L	10	U	1	U	1 U	1 U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	2 U	1 U	2 U	1 U	1 U	1 U		10 U	10 L	10	U	0.81	J	1 U	1 U
7440-39-3	BARIUM (DISSOLVED)	ug/l	47	39	32	19	34	110		40	40	42		44		38	45
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	1 U		1 U	1 L	1	U	1	U	2 U	1 U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	5.7		1 U	1 L	1	U	1	U	1 U	1 U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	35	1 U	7.3	1.9	43	1 U		5 U	5 L	5	U	25		22	3.5 J
7440-50-8	COPPER (DISSOLVED)	ug/l	NA	5 U	NA	NA	5 U	22	1	NA	NA	NA		NA		NA	NA
7439-92-1	LEAD (DISSOLVED)	ug/l	1 U	1 U	1 U	1 U	1 U	1 U		5 U	5 L	5	U	1	U	1 U	1 U
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	(0.2 U	0.2 L	0.2	U	0.2	J	0.71	0.2 U
7440-02-0	NICKEL (DISSOLVED)	ug/l	7.2	5.6	2.5	1.2	110	25		1.4 J	1.6 J	2.3	J	1.3		3.3	1.2
7782-49-2	SELENIUM (DISSOLVED)	ug/l	2 U	2 U	2 U	1 U	2 U	2 U	7	7.7 J	5.5 J	5.8	J	9.3	J	1 U	7.8
7440-22-4	SILER (DISSOLVED)	ug/l	1 U	0.1 U	1 U	1 U	0.1 U	0.1 U		5 U	5 L	1.1	J	1	U	1 U	1 U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	1 U	1 U	1 U	1 Ü	1 U	120		10 U	10 L	10	U	1	U	2 U	0.18 J
7440-66-6	ZINC (DISSOLVED)	ug/l	5 U	10 U	5 U	1.2 J	10 U	850		3.2 J	50 L	9.6	J	2.5	U	2.4 J	2.3 J
75-25-2	BROMOFORM	ug/l	NA	NA	NA NA	NA	NA	NA	0).51 J	1 1	1	U	NA		NA	NA
108-88-3	TOLUENE	ua/l	NA	NA	NA	NA	NA	NA		1 U	1 1	1	U	NA	1 1	NA	NA

Human Health and Enironmental Risk Characterization

	L	ocation:								DD-MW-203									DD-MW-204			
	Sa	mple ID:	DD-MW-201-R06	5-X	DD-MW-203-001	1-X	DD-MW-203-R02-X	DD-MW-203-R	03-X	DD-MW-203-R03-X-	FF_E	DD-MW-203-R04	4-X	DD-MW-203-R05-X		DD-MW-203-R06-X		DD-MW-204-001-X	DD-MW-204-R02-X	DD-MW-204-R05-)	х	DD-MW-205-001-X
	Samı	ple Date:	5/19/2008		6/26/2006		8/2/2006	8/14/2006		8/14/2006		6/5/2007		12/11/2007		5/19/2008		6/26/2006	6/5/2007	5/19/2008		6/26/2006
		Depth:	4 - 14 feet		8.5 - 8.5 feet		8.5 - 8.5 feet	10.5 - 10.5 fe	et	10.5 - 10.5 fee	t	4.5 - 12.5 feet	:	4.5 - 12.5 feet		4.5 - 12.5 feet		10.5 - 10.5 feet	7.5 - 13.5 feet	7.5 - 13.5 feet		9 - 9 feet
CAS Number	Analyte	Únits	Result	Q	Result	Q	Result Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result Q	Result C	Result	Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	5.2	U	10	UJ	NA	5.8	UJ	NA		5.2	U	5.1 L	JJ	5.1	U	10 U	5.1 L	0.84	J	10 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.31	U	0.94		NA	0.35	u	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
50-32-8	BENZO(A)PYRENE	ug/l	0.21	U	0.79		NA	0.23	u	0.23	u	0.2	U	0.2	U	0.2	U	0.21 U	0.2 L	0.2	U	0.2 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.31	U	0.74		NA	0.35	U	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.52	U	0.79		NA	0.58	U	0.57	u	0.4	U	0.51 U	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.31	U	0.76		NA	0.35	U	0.34	u	0.3	U	0.3	U	0.3	U	0.32 U	0.3 L	0.3	U	0.3 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.48	J	10	UJ	NA	5.8	U	NA		5.2	J	5.1	J	0.43	J	10 U	5.1	0.45	٦	10 U
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	U	5.1 l	U	5.1	U	10 U	5.1 L	5.1	U	10 U
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 I	U	100	U	110 U	100 L	100	U	100 U
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 I	U	100	U	110 U	100 L	100	U	100 U
PH912	C9-C18 ALIPHATICS	ug/l	100	U	100	U	NA	120	U	110	u	100	U	100 l	U	100	U	110 U	100 L	100	U	100 U
218-01-9	CHRYSENE	ug/l	1	U	1	U	NA	1.2	U	1.1	u	1	U	1 1	U	1	U	1.1 U	1 L	1	U	1 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.52	U	0.82		NA	0.58	U	0.57	u	0.4	U	0.51 l	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	J	5.1	J	5.1	U	10 U	5.1	5.1	U	10 U
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	5.2	U	10	UJ	NA	5.8	U	NA		5.2	U	5.1 l	U	5.1	U	10 U	5.1 L	5.1	U	10 U
206-44-0	FLUORANTHENE	ug/l	1	U	1	U	NA	1.2	U	1.1	u	1	U	1 1	U	1	U	1.1 U	1 L	1	U	1 U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.52	U	0.81		NA	0.58	U	0.57	u	0.4	U	0.51 l	U	0.51	U	0.53 U	0.4 L	0.51	U	0.51 U
85-01-8	PHENANTHRENE	ug/l	0.21	U	0.23		NA	0.23	U	0.23	u	0.2	U	0.2	U	0.2	U	0.21 U	0.2 L	0.2	U	0.2 U
129-00-0	PYRENE	ug/l	5.2	U	1	U	NA	1.2	U	1.1	u	1	U	5.1 l	U	5.1	U	1.1 U	1 L	5.1	U	1 U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	1	U	10	U	NA	10	U	NA		1	U	1 1	U	1	U	10 U	1 L	1	U	10 U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	1	U	10	U	NA	3.9	J	NA		1	U	1 1	U	1	U	10 U	1 L	2	U	10 U
7440-39-3	BARIUM (DISSOLVED)	ug/l	46		14		NA	13		NA		12		11		11		26	17	20		7.3 J
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	NA	0.096	J	NA		1	U	2 1	U	1	U	0.16 J	1 L	1	U	1 U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1	U	1	U	NA	1	U	NA		1	U	1 1	U	1	U	1 U	1 L	1	U	1 U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	2.1	J	5	U	NA	5	U	NA		16		15		0.63	J	5 U	11	1	U	5 U
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA	NA		NA		NA		NA		NA		NA	NA	NA		NA
7439-92-1	LEAD (DISSOLVED)	ug/l	1	U	5	U	NA	5	U	NA		1	U	1 1	U	1	U	5 U	1 L	1	U	5 U
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	U	NA	0.2	U	NA		0.2	U	0.2	U	0.2	U	0.2 U	0.2 L	0.2	Ū	0.2 U
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.3		1.6	J	NA	1.3	J	NA		1.1		2	U	1.2		1.9 J	1.1	1.2		1.3 J
7782-49-2	SELENIUM (DISSOLVED)	ug/l	6.8		10	Ü	NA	7.3	J	NA		1	U	1 1	Ū	1	U	10 U	1 L	2	U	10 U
7440-22-4	SILER (DISSOLVED)	ug/l	1	U	5	U	NA	1.4	J	NA		1	U	1 1	U	1	U	5 U	1 L	1	U	5 U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	1	U	10	U	NA	10	U	NA		1	U	2	U	0.25	J	10 U	1 L	1	Ū	10 U
7440-66-6	ZINC (DISSOLVED)	ug/I	2.7		50	U	NA	1.4	J	NA		2.4	J	1.9	J	5.3		1.6 J	2.1	5	U	50 U
75-25-2	BROMOFORM	ug/l	NA		1	U	NA	1	U	NA		NA		NA		NA		0.67 J	NA	NA		1 U
108-88-3	TOLUENE	ua/l	NA		1	U	NA	1	U	NA		NA		NA		NA		1 U	NA	NA		1 U

Human Health and Enironmental Risk Characterization

	L	ocation:	D	D-MV	V-205									DD-MW-206									
	Sa	mple ID:	DD-MW-205-R02-	-X	DD-MW-205-R0)3-X	DD-MW-205-R05-X	DD-MW-206-00	1-X	DD-MW-206-R02	2-D	DD-MW-206-R02	2-X	DD-MW-206-R03->	(DD-MW-206-R03-X-I	D	DD-MW-206-R04-X		DD-MW-206-R05-X	DD-MW-207-001-X	K	DD-MW-207-R02-X
	Samp	ole Date:	8/14/2006		6/6/2007		5/19/2008	6/26/2006		8/2/2006		8/2/2006		8/14/2006		8/14/2006		6/5/2007		5/19/2008	6/27/2006		8/2/2006
		Depth:	10 - 10 feet		5 - 11 feet		5 - 11 feet	8 - 8 feet		10 - 10 feet		10 - 10 feet		10 - 10 feet		10 - 10 feet		5 - 11 feet		5 - 11 feet	10 - 10 feet		12.5 - 12.5 feet
CAS Number	Analyte	Units	Result	Q	Result	Q	Result Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result Q	2	Result Q	Result	Q	Result Q
51-28-5	2,4-DINITROPHENOL	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	UJ	5.1 l	UJ	5.2 U	J	5.1 U	10	U	NA
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
50-32-8	BENZO(A)PYRENE	ug/l	0.2	U	0.2	U	0.2 U	0.21	U	NA		NA		0.2	U	0.2	U	0.21 U	J	0.2 U	0.2	U	NA
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.3 U	0.32	U	NA		NA		0.3	U	0.3	U	0.31 U	J	0.3 U	0.3	U	NA
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	10	U	5.1	J	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 J	J	0.49 J	10	U	NA
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 U	J	5.1 U	10	U	NA
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
PH912	C9-C18 ALIPHATICS	ug/l	100	U	100	U	100 U	110	U	NA		NA		100	U	100	U	100 U	J	100 U	100	U	NA
218-01-9	CHRYSENE	ug/l	1	U	1	U	1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	10	U	5.1	J	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 J	J	5.1 U	10	U	NA
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	10	U	5.1	U	5.1 U	11	UJ	NA		NA		5.1	U	5.1	U	5.2 U	J	5.1 U	10	U	NA
206-44-0	FLUORANTHENE	ug/l	1	U	1	U	1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.51	U	0.4	U	0.51 U	0.53	U	NA		NA		0.51	U	0.51	U	0.41 U	J	0.51 U	0.51	U	NA
85-01-8	PHENANTHRENE	ug/l	0.2	U	0.2	U	0.2 U	0.21	U	NA		NA		0.2	U	0.2	U	0.21 U	J	0.2 U	0.2	U	NA
129-00-0	PYRENE	ug/l	1	U	1	U	5.1 U	1.1	U	NA		NA		1	U	1	U	1 U	J	5.1 U	1	U	NA
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	10	U	1	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	10	U	NA
7440-38-2	ARSENIC (DISSOLVED)	ug/l	10	U	2	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	4.7	J	NA
7440-39-3	BARIUM (DISSOLVED)	ug/l	6.1	J	18		7.6	23		NA		NA		27		28		26		16	38		NA
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	1 U	1	U	NA		NA		1	U	1	U	1 U	J	1 U	1	U	NA
7440-43-9	CADMIUM (DISSOLVED)	ug/l	1	U	1	U	1 U	1	U	NA		NA		1	U	1	U	1 U	J	1 U	0.69	J	NA
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	5	U	8.4		0.52 J	5	U	NA		NA		5	U	5	U	23		0.51 J	5	U	NA
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA	NA		NA		NA		NA		NA		NA		NA	NA		NA
7439-92-1	LEAD (DISSOLVED)	ug/l	5	U	1	U	1 U	5	U	NA		NA		5	U	5	U	1 U	J	1 U	5	U	NA
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	U	0.2 U	0.2	U	NA		NA		0.2	U	0.2	U	0.2 U	J	0.2 U	0.2	U	NA
7440-02-0	NICKEL (DISSOLVED)	ug/l	1.3	J	1.4		0.59 J	1.4	J	NA		NA		1.3	J	10	U	0.91 J	J	0.69 J	4.7	J	NA
7782-49-2	SELENIUM (DISSOLVED)	ug/l	10	U	2	U	1 U	10	U	NA		NA		10	U	10	U	1 U	J	1 U	10	U	NA
7440-22-4	SILER (DISSOLVED)	ug/l	5	U	1	U	1 U	5	U	NA		NA		5	U	1.5	J	1 U	J	1 U	5	U	NA
7440-62-2	ANADIUM (DISSOLVED)	ug/l	10	Ū	1	U	0.29 J	10	U	NA		NA		10	U	10	Ü	1 U	J	1 U	10	Ū	NA
7440-66-6	ZINC (DISSOLVED)	ug/l	50	U	5	U	1.6 J	4.2	J	NA		NA		1.8	J	9.3	J	2.7		1.9 J	3.5	J	NA
75-25-2	BROMOFORM	ug/l	1	Ū	NA		NA	0.83	J	NA		NA		1	U	1	U	NA		NA	1	Ü	NA
108-88-3	TOLUENE	ua/l	1	U	NA		NA	1	U	NA		NA		1	U	0.52	J	NA		NA	1	U	NA

Human Health and Enironmental Risk Characterization

		Location:			DD-MW-207												DD-MW-208-R01-	001-D				
		Sample ID:	DD-MW-207-R	03-X	DD-MW-207-R04	I-X	DD-MW-207-R05	5-D	DD-MW-207-R05	5-X	DD-MW-207-R0)6-X	DD-MW-208-R01-0	01-D	DD-MW-208-R01-0	01-X	DD-MW-208-R0	2-X	DD-MW-208-R0	03-X	DD-MW-208-R0	4-X
	Sa	ample Date:	8/14/2006		6/6/2007		12/11/2007		12/11/2007		5/19/2008		6/25/2007		6/25/2007		7/23/2007		12/11/2007		5/19/2008	
		Depth:	12.5 - 12.5 fe	eet	5 - 15 feet		5 - 15 feet		5 - 15 feet		5 - 15 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet		3 - 13 feet	
CAS Number	Analyte	Units	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q	Result	Q
51-28-5	2,4-DINITROPHENOL	ug/l	10	U	5.1	U	5.5	U	5.1	UJ	5.2	U	5.1	U	5.1	UJ	5.2	U	5.1	UJ	5.1	U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.3	U	0.22	J	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
50-32-8	BENZO(A)PYRENE	ug/l	0.2	U	0.2	U	0.22	U	0.2	U	0.21	U	0.2	U	0.2	U	0.21	U	0.2	U	0.2	U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3	U	0.32	U	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.51	U	0.42	U	0.55	U	0.51	U	0.52	U	0.4	U	0.41	U	0.21	J	0.51	U	0.51	U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.3	U	0.3	U	0.33	U	0.3	U	0.31	U	0.3	U	0.3	U	0.31	U	0.3	U	0.3	U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	10	U	5.1	J	5.5	J	5.1	J	0.51	J	5.1	U	5.1	U	0.78	J	5.1	J	1.6	J
85-68-7	BUTYLBENZYLPHTHALATE	ug/l	10	U	5.1	U	5.5	U	5.1	U	5.2	U	5.1	U	5.1	U	0.34	J	5.1	U	5.1	U
EPH1122	C11-C22 AROMATICS, ADJUSTED	ug/l	100	U	110	U	100	U	100	U	100	U	290		460		100	U	NA		100	U
EPH1936	C19-C36 ALIPHATICS	ug/l	100	U	110	U	100	U	100	U	100	U	100	U	130		100	U	100	U	100	U
PH912	C9-C18 ALIPHATICS	ug/l	100	U	110	U	100	U	100	U	100	U	100	U	100	U	100	U	100	U	100	U
218-01-9	CHRYSENE	ug/l	1	U	0.24	J	1.1	U	1	U	1	U	1	U	1	U	1	U	1	U	1	U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.51	U	0.42	U	0.55	U	0.51	U	0.52	U	0.4	U	0.41	U	0.42	J	0.51	U	0.51	U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	10	U	5.1	J	5.5	J	5.1	J	5.2	U	5.1	J	5.1	J	5.2	J	5.1	J	1.5	J
117-84-0	DI-N-OCTYL PHTHALATE	ug/l	10	U	0.39	J	5.5	U	5.1	U	5.2	U	5.1	U	5.1	U	5.2	U	5.1	U	5.1	U
206-44-0	FLUORANTHENE	ug/l	1	U	0.28	J	1.1	U	1	U	1	U	1.885	J	1.875	J	1	U	1	U	1	U
193-39-5	INDENO(1,2,3-CD)PYRENE	ug/l	0.51	U	0.42	U	0.55	U	0.51	U	0.52	U	0.4	U	0.41	U	0.38	J	0.51	U	0.51	U
85-01-8	PHENANTHRENE	ug/l	0.2	U	0.2	U	0.22	U	0.2	U	0.21	U	0.62	J	0.37	J	0.21	U	0.2	U	0.2	U
129-00-0	PYRENE	ug/l	1	U	0.29	J	5.5	U	5.1	U	5.2	U	0.28	J	1	U	1	U	5.1	U	5.1	U
7440-36-0	ANTIMONY (DISSOLVED)	ug/l	2.3	U	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-38-2	ARSENIC (DISSOLVED)	ug/l	7.7	J	5.7		10		10		3.9		1.7	J	1.7	J	NA		1	U	1	U
7440-39-3	BARIUM (DISSOLVED)	ug/l	44		38		44		45		31		27		26		NA		25		31	
7440-41-7	BERYLLIUM (DISSOLVED)	ug/l	1	U	1	U	2	U	2	U	1	U	2	U	1	U	NA		2	U	1	U
7440-43-9	CADMIUM (DISSOLVED)	ug/l	0.84	J	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-47-3	CHROMIUM (DISSOLVED)	ug/l	5	U	4.9		9.1		10		1.3		9.8		11		NA		25		0.36	J
7440-50-8	COPPER (DISSOLVED)	ug/l	NA		NA		NA		NA		NA		NA		NA		NA		NA		NA	
7439-92-1	LEAD (DISSOLVED)	ug/l	5	U	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	0.46	J
7439-97-6	MERCURY (DISSOLVED)	ug/l	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	NA		0.2	U	0.2	U
7440-02-0	NICKEL (DISSOLVED)	ug/l	4	J	1	U	2	U	2	U	0.43	J	3.6		3.2		NA		0.96	J	2	
7782-49-2	SELENIUM (DISSOLVED)	ug/l	10	U	1	U	1	U	1	U	1	U	2	U	2	U	NA		1	U	1	U
7440-22-4	SILER (DISSOLVED)	ug/l	5	U	1	U	1	U	1	U	1	U	1	U	1	U	NA		1	U	1	U
7440-62-2	ANADIUM (DISSOLVED)	ug/l	0.92	J	1	U	2	U	2	U	1	U	0.91	J	1	J	NA		2	U	0.52	J
7440-66-6	ZINC (DISSOLVED)	ug/l	3.5	J	2.5	U	1.2	J	1.5	J	1.9	J	9.4		7.7		NA		20		110	
75-25-2	BROMOFORM	ug/l	1	U	NA		NA		NA		NA		5	U	5	U	NA		NA		NA	
108-88-3	TOLUENE	ug/l	1	U	NA		NA		NA		NA		5	U	5	U	NA		NA		NA	T

Attachment A(3)
Analytical Data - DDA Earthworm
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	Lo	cation	BP-BO-007	,	DD-BO-001		DD-BO-002	2	BP-BO-00)7
	Sam	ple ID	BP-BO-007-00	1-X	DD-BO-001-00)1-X	DD-BO-002-00)1-X	SH-BO-001-0)01-X
	Sample	e Date	9/22/2006		9/21/2006		9/22/2006		9/22/200	6
		Depth			0 - 1 feet		0 - 1 feet			
Cas Number	Analyte	Units	Results	Q	Results	Q	Results	Q	Results	Q
7440-36-0	ANTIMONY	mg/kg			0.26	В	0.10	В	0.091	В
7440-38-2	ARSENIC	mg/kg			1.30	V	0.82	V	0.34	V
7440-39-3	BARIUM	mg/kg							6.60	В
7440-41-7	BERYLLIUM	mg/kg			0.24	В	0.14	В	0.041	В
7440-43-9	CADMIUM	mg/kg			8.10	V	3.80	V	1.30	V
7440-47-3	CHROMIUM	mg/kg			16.20	V	10.30	V	4.100	V
7440-50-8	COPPER	mg/kg			14.80	V	3.90	V	2.00	V
7439-92-1	LEAD	mg/kg			36.20	V	101	V	28.30	V
7440-02-0	NICKEL	mg/kg			9.70	V	6.90	V	2.60	V
7782-49-2	SELENIUM	mg/kg							0.91	V
7440-22-4	SILVER	mg/kg			0.18	V	0.040	В	0.058	В
7440-28-0	THALLIUM	mg/kg			0.024	В	0.024	В	0.025	В
7440-62-2	VANADIUM	mg/kg			5.20	V	5.00	V	3.00	V
7440-66-6	ZINC	mg/kg			168	V	114	V	81.20	V
7439-97-6	MERCURY	mg/kg			0.45	V	0.10	V	0.066	V
% SOLIDS	% SOLIDS	%			37.30	V	38.90	V	28.40	V
	%LIPIDS DETERMINATION	%	1.60	٧	0.90	V	1.00	V		

Notes:

mg/kg - milligrams per kilogram

Q - qualifier

DDA - Demoilition Debris Area

1 of 1 11/2/2011



Background Arsenic Evaluation

Identification of Background Wells

Background is defined as those levels of oil and hazardous material (arsenic in this case) that would exist in the absence of the disposal site of concern, also usually described as naturally occurring or anthropogenic levels of "contaminants". Arsenic is naturally occurring in soil and groundwater throughout Massachusetts as a result of geologic processes. Monitoring Wells that appear to represent groundwater not influenced by the Site include the following locations:

- CP-MW-101 is located adjacent to the north side of Ruckaduck Pond, and upgradient of known manufacturing and disposal areas on the Site;
- DD-MW-204 and -205 are located upgradient of the DDA;
- MB-MW-360 and -361 are located east of the Neponset River; and
- MB-MW-368 is located near the southeast side of Ruckaduck Pond, upgradient from the former manufacturing buildings and crossgradient from the South Rail Spur source area.

Evidence of no Site impacts for these wells includes no organic contaminants detected, and positions that are upgradient or crossgradient of known source areas based on groundwater flow directions mapped in the Phase II and Phase III reports. All of these wells are screened in the shallow sandy aquifer where other Site monitoring wells are screened. The average arsenic concentration for each background well is listed in Table B-1.

Methods and Results

EPA's ProUCL software (version 4.1) was used to calculate upper tolerance levels (UTLs) for groundwater collected between 1992 and 2009. The UTL calculated represents the value below which 90% of background values are expected to fall with a 95% confidence. All UTLs provided by ProUCL for each distribution were evaluated. Note that ProUCL guidance recommends the use of either the largest value or the second largest value for a UTL, while cautioning that the largest UTL may result in an overestimate. For the purposes of this evaluation, the second largest recommended UTL was used to represent background concentrations at this site.

Results from ProUCL are provided in Table B-2 and indicate that groundwater data follow a normal, gamma and lognormal distribution (likely due to the relatively small sample size of the dataset; represented by 6 observations). Parametric statistics predict background concentrations at 0.0143 mg/l, 0.049 mg/l and 0.102 mg/l assuming a normal, gamma, and lognormal distribution, respectively. Since the dataset size is limited an alternate and more appropriate approach maybe to use non-parametric statistics to estimate background concentrations. The recommended UTL from this set of statistical tests calculates a background arsenic concentration of 0.0159 mg/l.

Table B-1
Background Results of Arsenic
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

Sample ID	Units	Results	Qualifier
MW-101	mg/L	0.0103	
MW-204	mg/L	0.001	U
MW-205	mg/L	0.001	U
MW-360	mg/L	0.0039	
MW-361	mg/L	0.00079	
MW-368	mg/L	0.0086	

mg/L - milligrams per liter U - Value is non-detect

Table B-2. General Background Statistics for Data Sets with Non-Detects

User Selected Options

ProUCL.wst From File Full Precision OFF Confidence Coefficient 95% 90% Coverage Different or Future K Values Number of Bootstrap Operations 2000

Arsenic

General Statistics

Number of Valid Data 6 Number of Detected Data 4 Number of Distinct Detected Data 4 Number of Non-Detect Data 2 Tolerance Factor 3.006 Percent Non-Detects 33 33%

Raw Statistics

Minimum Detected 0.00079 Maximum Detected 0.0103 Mean of Detected 0.0059 SD of Detected 0.00435 Minimum Non-Detect 0.001 Maximum Non-Detect 0.001

Log-transformed Statistics

Minimum Detected -7.143 Maximum Detected -4.576 Mean of Detected -5.505 SD of Detected 1.171 Minimum Non-Detect -6.908 Maximum Non-Detect -6.908

Warning: There are only 4 Distinct Detected Values in this data Note: It should be noted that even though bootstrap may be performed on this data set the resulting calculations may not be reliable enough to draw conclusions

It is recommended to have 10-15 or more distinct observations for accurate and meaningful results.

Background Statistics

Normal Distribution Test with Detected Values Only Shapiro Wilk Test Statistic 0.945

5% Shapiro Wilk Critical Value 0.748

Data appear Normal at 5% Significance Level

Lognormal Distribution Test with Detected Values Only Shapiro Wilk Test Statistic 0.876

5% Shapiro Wilk Critical Value 0.748

Data appear Lognormal at 5% Significance Level

Assuming Normal Distribution

DL/2 Substitution Method Mean 0.0041 SD 0.00437 95% UTL 90% Coverage 0.0172 95% UPL (t) 0.0136 90% Percentile (z) 0.0097 95% Percentile (z) 0.0113 99% Percentile (z) 0.0143

Maximum Likelihood Estimate(MLE) Method

Mean 0.00765

95% UTL with 90% Coverage 0.0158

SD 0.00271

95% UPL (t) 0.0135 90% Percentile (z) 0.0111 95% Percentile (z) 0.0121

99% Percentile (z) 0.0139

Gamma Distribution Test with Detected Values Only

k star (bias corrected) 0.539 Theta Star 0.0109 nu star 4.309

A-D Test Statistic 0.343 5% A-D Critical Value 0.662 K-S Test Statistic 0.276

5% K-S Critical Value 0.4 Data appear Gamma Distributed at 5% Significance Level

Assuming Gamma Distribution

Gamma ROS Statistics with Extrapolated Data

Mean 0.00412 Median 0.00252 SD 0.00436 k star 0.315 Theta star 0.0131 Nu star 3.778

95% Percentile of Chisquare (2k) 2.836

90% Percentile 0.0121 95% Percentile 0.0186 99% Percentile 0.0353

Assuming Lognormal Distribution

DL/2 Substitution Method Mean (Log Scale) -6.204 SD (Log Scale) 1.412 95% UTL 90% Coverage 0.141 95% UPL (t) 0.0437 90% Percentile (z) 0.0123 95% Percentile (z) 0.0206 99% Percentile (z) 0.054

> Log ROS Method Mean in Original Scale 0.00423 SD in Original Scale 0.00425

95% UTL with 90% Coverage 0.102 95% BCA UTL with 90% Coverage 0.0103

95% Bootstrap (%) UTL with 90% Coverage 0.0103 95% UPL (t) 0.0362 90% Percentile (z) 0.0118 95% Percentile (z) 0.0186 99% Percentile (z) 0.0436

Data Distribution Test with Detected Values Only

Data appear Normal at 5% Significance Level

Nonparametric Statistics

Kaplan-Meier (KM) Method Mean 0.0042 SD 0.00391

SE of Mean 0.00184

95% KM UTL with 90% Coverage 0.0159 95% KM Chebyshev UPL 0.0226 95% KM UPL (t) 0.0127

90% Percentile (z) 0.0092 95% Percentile (z) 0.0106 99% Percentile (z) 0.0133

Gamma ROS Limits with Extrapolated Data

95% Wilson Hilferty (WH) Approx. Gamma UPL 0.0273 95% Hawkins Wixley (HW) Approx. Gamma UPL 0.0378

95% WH Approx. Gamma UTL with 90% Coverage 0.049 95% HW Approx. Gamma UTL with 90% Coverage 0.0784



Attachment C TCDD TEQ Calculations Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

		Location			DD-GP-003	DD-GP-004	DD-GP-006	DD-GP-009	DD-GP-207	DD-SB-206	DD-SS-003	DD-SS-005
		Sample ID			DDA-GP-3 RR & Dup	DDA-GP-4	DDA-GP-6	DDA-GP-9 RR & Dup	DD-GP-207-001-X	DD-SB-206-003-X	DD-SS-001-001-X	DD-SS-005-001-X
		Sample Date			5/16/2005	5/16/2005	5/16/2005	5/16/2005	6/9/2006	6/19/2006	12/8/2005	9/21/2006
CAS		Depth	Mammalian	Avian	0 - 3 feet	2 - 4 feet	0 - 3 feet	0 - 3 feet	0 - 2 feet	0 - 2 feet	0 - 3 feet	0 - 0.5 feet
Number	Analyte	units	TEF	TEF	Results	Results	Results	Results	Results	Results	Results	Results
								<u> </u>				
1746-01-6	2,3,7,8-TCDD	mg/kg	1	1	2.35E-07 U	2.50E-07 U	1.30E-07 U	1.15E-07 U	8.50E-08 U	5.50E-06	NA	NA
40321-76-4	1,2,3,7,8-PeCDD	mg/kg	1	1	3.50E-06 J	5.00E-07 U	3.35E-07 U	1.35E-07 U	3.85E-07 U	1.70E-05	NA	NA
39227-28-6	1,2,3,4,7,8-HxCDD	mg/kg	0.1	0.05	1.35E-06 U	3.05E-07 U	2.60E-07 U	1.15E-07 U	2.20E-07 U	7.30E-06 J	NA	NA
57653-85-7	1,2,3,6,7,8-HxCDD	mg/kg	0.1	0.01	1.40E-05 V	3.20E-07 U	1.40E-06 U	2.55E-07 U	2.00E-07 U	4.00E-05	NA	NA
19408-74-3	1,2,3,7,8,9-HxCDD	mg/kg	0.1	0.01	6.00E-06 V	3.20E-07 U	9.00E-07 U	2.30E-07 U	2.00E-07 U	2.40E-05	NA	NA
35822-46-9	1,2,3,4,6,7,8-HpCDD	mg/kg	0.01	0.001	3.10E-04 V	5.70E-06 J	7.10E-05 V	6.30E-06 V	9.20E-06 v	8.30E-04	NA	NA
3268-87-9	OCDD	mg/kg	0.0003	0.0001	2.70E-03 V	4.20E-05 V	1.10E-03 V	4.40E-05 V	9.80E-05 v	7.20E-03	NA	NA
51207-31-9	2,3,7,8-TCDF	mg/kg	0.1	1	2.10E-06	6.10E-07	1.30E-06	2.90E-07	4.90E-07	3.70E-05	NA	NA
57117-41-6	1,2,3,7,8-PeCDF	mg/kg	0.03	0.01	9.50E-07 U	2.55E-07 U	7.50E-07 U	1.15E-07 U	2.70E-07 U	1.60E-05	NA	NA
57117-31-4	2,3,4,7,8-PeCDF	mg/kg	0.3	1	5.70E-06 J	2.55E-07 U	3.80E-06 J	3.25E-07 U	2.70E-07 U	7.10E-05	NA	NA
70648-26-9	1,2,3,4,7,8-HxCDF	mg/kg	0.1	0.1	7.20E-06 V	9.00E-07 U	4.30E-06 J	6.50E-07 U	9.50E-07 U	2.70E-05	NA	NA
57117-44-9	1,2,3,6,7,8-HxCDF	mg/kg	0.1	0.1	8.90E-06 V	3.50E-07 U	4.20E-06 J	3.80E-07 U	3.40E-07 U	9.00E-05	NA	NA
72918-21-9	1,2,3,7,8,9-HxCDF	mg/kg	0.1	0.1	4.55E-07 U	3.20E-07 U	5.50E-07 U	2.40E-07 U	1.80E-07 U	3.60E-06 U	NA	NA
60851-34-5	2,3,4,6,7,8-HxCDF	mg/kg	0.1	0.1	8.70E-06 V	5.50E-07 U	5.50E-06 J	4.15E-07 U	6.50E-07 U	1.10E-04	NA	NA
67562-39-4	1,2,3,4,6,7,8-HpCDF	mg/kg	0.01	0.01	9.00E-05 V	6.50E-06 J	2.30E-05 V	6.30E-06 V	6.80E-06 v	3.10E-04	NA	NA
55673-89-7	1,2,3,4,7,8,9-HpCDF	mg/kg	0.01	0.01	4.00E-06 J	2.30E-07 U	5.50E-07 U	1.85E-07 U	2.50E-07 U	1.20E-05	NA	NA
39001-02-0	OCDF	mg/kg	0.0003	0.0001	4.60E-04 V	3.90E-06 U	2.50E-05 V	6.40E-06 J	2.00E-05 v	1.50E-03	NA	NA
							•		•	•	•	•
	Mammalian TEQ	mg/kg			0.000015	0.000013	0.000048	0.0000075	0.0000011	0.00009230	NA	NA
	•							•		•		
	Avian TEQ	mg/kg			0.000016	0.000019	0.000075	0.00000112	0.000016	0.00015965	NA	NA

mg/kg-milligrams per kilogram
TEF - Toxicity equivalency factor
TEQ - Toxic equivalency quotient

NA - not analyzed
(1) TEFs presented are the World Health Organization
(WHO) 2005 values from Van den Berg et al., 2006

(2) Mammailian TEQs were calculated by multiplying the results for each individual congener by the applicable Mammalian TEFs and summing the results.

Avian TEQs were calculated using the same process and the Avian TEFs.

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Attachment C TCDD TEQ Calculations Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

		Location			DD-SS-007	DD-SS-012	DD-SS-014	DD-T	P-013	DD-TP-003	DD-TP-004	DD-TP-005
		Sample ID			DD-SS-007-001-X & Dup	DD-SS-012-001-X	DD-SS-014-001-X	DD-TP-001-001-X	DD-TP-001-002-X	DD-TP03-2S	DD-TP04-2S	DD-TP05-2S
		Sample Date			10/26/2006	5/30/2007	5/30/2007	11/21/2005	11/21/2005	12/20/2004	12/20/2004	12/20/2004
CAS		Depth	Mammalian	Avian	0 - 0.5 feet	0 - 2 feet	0 - 2 feet	0 - 2 feet	2 - 4 feet	0 - 3 feet	0 - 3 feet	0 - 3 feet
Number	Analyte	units	TEF	TEF	Results	Results	Results	Results	Results	Results	Results	Results
1746-01-6	2,3,7,8-TCDD	mg/kg	1	1	1.75E-06 v	1.55E-07 U	3.90E-06 v	NA	NA	5.90E-07 J	NA	4.20E-06 V
40321-76-4	1,2,3,7,8-PeCDD	mg/kg	1	1	6.90E-06 v	3.10E-07 U	1.80E-05 v	NA	NA	1.50E-06 J	NA	1.40E-05 V
39227-28-6	1,2,3,4,7,8-HxCDD	mg/kg	0.1	0.05	1.35E-06 U	2.45E-07 U	7.70E-06 J	NA	NA	1.10E-06 J	NA	1.30E-05 V
57653-85-7	1,2,3,6,7,8-HxCDD	mg/kg	0.1	0.01	1.25E-05 v	4.15E-07 U	3.60E-05 v	NA	NA	3.30E-06 J	NA	2.70E-05 V
19408-74-3	1,2,3,7,8,9-HxCDD	mg/kg	0.1	0.01	7.15E-06 v	3.15E-07 U	2.40E-05 v	NA	NA	3.20E-06 J	NA	4.00E-05 V
35822-46-9	1,2,3,4,6,7,8-HpCDD	mg/kg	0.01	0.001	1.50E-04 v	7.20E-06 v	1.80E-04 v	NA	NA	2.10E-05 B	NA	6.70E-04 B
3268-87-9	OCDD	mg/kg	0.0003	0.0001	1.25E-03 v	5.40E-05 v	1.40E-03 v	NA	NA	1.00E-04 B	NA	5.70E-03
51207-31-9	2,3,7,8-TCDF	mg/kg	0.1	1	1.30E-05	1.40E-06	2.50E-05	NA	NA	4.50E-06 V	NA	5.80E-05 V
57117-41-6	1,2,3,7,8-PeCDF	mg/kg	0.03	0.01	5.75E-06 J	4.70E-07 U	2.10E-05 v	NA	NA	3.30E-06 J	NA	3.30E-05 V
57117-31-4	2,3,4,7,8-PeCDF	mg/kg	0.3	1	2.30E-05 v	1.00E-06 U	4.70E-05 v	NA	NA	6.80E-06 V	NA	6.60E-05 V
70648-26-9	1,2,3,4,7,8-HxCDF	mg/kg	0.1	0.1	9.40E-06 v	8.50E-07 U	2.90E-05 v	NA	NA	2.00E-05 V	NA	1.60E-04 V
57117-44-9	1,2,3,6,7,8-HxCDF	mg/kg	0.1	0.1	2.55E-05 v	8.00E-07 U	4.70E-05 v	NA	NA	7.40E-06 V	NA	6.40E-05 V
72918-21-9	1,2,3,7,8,9-HxCDF	mg/kg	0.1	0.1	3.50E-07 U	3.15E-07 U	1.35E-06 U	NA	NA	8.50E-08 U	NA	2.90E-06 J
60851-34-5	2,3,4,6,7,8-HxCDF	mg/kg	0.1	0.1	3.15E-05 v	9.50E-07 U	5.30E-05 v	NA	NA	7.80E-06 V	NA	7.70E-05 V
67562-39-4	1,2,3,4,6,7,8-HpCDF	mg/kg	0.01	0.01	5.65E-05 v	3.50E-06 J	9.20E-05 v	NA	NA	6.20E-05 V	NA	2.70E-04 V
55673-89-7	1,2,3,4,7,8,9-HpCDF	mg/kg	0.01	0.01	4.80E-06 J	1.80E-07 U	6.70E-06 J	NA	NA	2.80E-06 J	NA	1.70E-05 V
39001-02-0	OCDF	mg/kg	0.0003	0.0001	1.05E-04 v	2.60E-06 U	1.30E-04 v	NA	NA	3.60E-05 B	NA	1.50E-04 B
					·	•	•		<u>.</u>		•	
	Mammalian TEQ	mg/kg			0.000028	0.0000014	0.000062	NA	NA	0.000099	NA	0.000095
	.									-		
	Avian TEQ	mg/kg			0.000053	0.0000032	0.000109	NA	NA	0.000018	NA	0.00018

mg/kg-milligrams per kilogram
TEF - Toxicity equivalency factor
TEQ - Toxic equivalency quotient

NA - not analyzed
(1) TEFs presented are the World Health Organization
(WHO) 2005 values from Van den Berg et al., 2006

(2) Mammailian TEQs were calculated by multiplying the results for each individual congener by the applicable Mammalian TEFs and summing the results.

Avian TEQs were calculated using the same process and the Avian TEFs.

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Attachment C TCDD TEQ Calculations Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

		Location			DD-TP-007	DD-TP-008	DD-TP-009	DD-TP-012	DD-TP-201	DD-T	P-202
		Sample ID			DD-TP07-2S-D & Dup	DD-TP08-2S	DD-TP09-2S-D & Dup	DD-TP12-5W	DD-TP-201-001-X	DD-TP-202-002-X	DD-TP-202-003-X
		Sample Date			12/20/2004	12/20/2004	12/20/2004	12/20/2004	6/6/2006	6/28/2006	6/28/2006
CAS		Depth	Mammalian	Avian	0 - 3 feet	0 - 3 feet	0 - 3 feet	0 - 3 feet	1 - 2 feet	2 - 3 feet	2 - 3 feet
Number	Analyte	units	TEF	TEF	Results	Results	Results	Results	Results	Results	Results
1746-01-6	2,3,7,8-TCDD	mg/kg	1	1	6.70E-07 J	NA	NA	NA	NA	NA	NA
40321-76-4	1,2,3,7,8-PeCDD	mg/kg	1	1	3.40E-06 J	NA	NA	NA	NA	NA	NA
39227-28-6	1,2,3,4,7,8-HxCDD	mg/kg	0.1	0.05	2.10E-06 J	NA	NA	NA	NA	NA	NA
57653-85-7	1,2,3,6,7,8-HxCDD	mg/kg	0.1	0.01	1.50E-05 V	NA	NA	NA	NA	NA	NA
19408-74-3	1,2,3,7,8,9-HxCDD	mg/kg	0.1	0.01	7.40E-06 V	NA	NA	NA	NA	NA	NA
35822-46-9	1,2,3,4,6,7,8-HpCDD	mg/kg	0.01	0.001	3.10E-04 B	NA	NA	NA	NA	NA	NA
3268-87-9	OCDD	mg/kg	0.0003	0.0001	7.60E-03	NA	NA	NA	NA	NA	NA
51207-31-9	2,3,7,8-TCDF	mg/kg	0.1	1	5.90E-06 V	NA	NA	NA	NA	NA	NA
57117-41-6	1,2,3,7,8-PeCDF	mg/kg	0.03	0.01	3.30E-06 J	NA	NA	NA	NA	NA	NA
57117-31-4	2,3,4,7,8-PeCDF	mg/kg	0.3	1	8.30E-06 V	NA	NA	NA	NA	NA	NA
70648-26-9	1,2,3,4,7,8-HxCDF	mg/kg	0.1	0.1	1.70E-05 V	NA	NA	NA	NA	NA	NA
57117-44-9	1,2,3,6,7,8-HxCDF	mg/kg	0.1	0.1	8.50E-06 V	NA	NA	NA	NA	NA	NA
72918-21-9	1,2,3,7,8,9-HxCDF	mg/kg	0.1	0.1	1.20E-07 U	NA	NA	NA	NA	NA	NA
60851-34-5	2,3,4,6,7,8-HxCDF	mg/kg	0.1	0.1	8.80E-06 V	NA	NA	NA	NA	NA	NA
67562-39-4	1,2,3,4,6,7,8-HpCDF	mg/kg	0.01	0.01	6.30E-05 V	NA	NA	NA	NA	NA	NA
55673-89-7	1,2,3,4,7,8,9-HpCDF	mg/kg	0.01	0.01	4.20E-06 J	NA	NA	NA	NA	NA	NA
39001-02-0	OCDF	mg/kg	0.0003	0.0001	1.70E-04 B	NA	NA	NA	NA	NA	NA
	Mammalian TEQ	mg/kg			0.000019	NA	NA	NA	NA	NA	NA
	Avian TEQ	mg/kg			0.000024	NA	NA	NA	NA	NA	NA

mg/kg-milligrams per kilogram
TEF - Toxicity equivalency factor
TEQ - Toxic equivalency quotient

NA - not analyzed
(1) TEFs presented are the World Health Organization
(WHO) 2005 values from Van den Berg et al., 2006

(2) Mammailian TEQs were calculated by multiplying the results for each individual congener by the applicable Mammalian TEFs and summing the results.

Avian TEQs were calculated using the same process and the Avian TEFs.

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Attachment D
Exposure and Risk Estimates Associated With Soil Contact
Bird Machine Company - DDA
100 Neponset St, Walpole MA
Trespasser Older Child
Surface Soil
DDA current
0-3 feet

Receptor:	Trespasser Older Child	▼
Medium:	Surface Soil	•
Exposure Area:	DDA current	▼
Depth:	0-3 feet	<u></u>
Duration:	Chronic	▼

Parameter	Definition	Units	Value	Comment	
IRsoil	Soil Ingestion Rate	mg/d	50		
SA	Soil Dermal Contact Skin Exposed	cm2/d	4260		
AF	Soil Dermal Contact Adherence Rate	mg/cm2	0.14		
EF	Soil Exposure Frequency	d/y	50		
EP	Soil Exposure Period - Cancer	У	7		
EP	Soil Exposure Period - Non-Cancer	y	7		
ATc	Soil Averaging Time - Cancer	d	25550		
ATn	Soil Averaging Time - Non-Cancer	d	2555		
BW	Body Weight	kg	39.9		
CF	Conversion Factor	kg/mg	0.000001		

$$ADD_{ing} = \frac{C_{soil} \times CF \times IR_{soil} \times RAF_{os} \times EF \times EP}{AT \times BW}$$

$$ADD_{der} = \frac{C_{soil} \times CF \times SA \times AF \times RAF_{ds} \times EF \times EP}{AT \times BW}$$

$$HI_{ing} = \frac{ADD_{ing}}{RfD}$$

$$HI_{der} = \frac{ADD_{der}}{RfD}$$

$$HI = HI_{ing} + HI_{der}$$

$$Risk_{ing} = ADD_{ing} \times CSF$$

$$Risk_{der} = ADD_{der} \times CSF$$

$$Risk_{er} = Risk_{ing} + Risk_{der}$$

	EPC			Incidental In	gestion					Dermal Co	ontact					Total	
	Surface Soil																
	DDA current																
	0-3 feet	RfD	CSF	RAFosc	ADDing-c	Risking	RAFosnc	ADDing-nc	Hling	RAFdsc	ADDder-c	Riskder	RAFdsnc	ADDder-nc	Hlder	Risk (Soil)	HI (Soil)
Compound	(mg/kg)	(mg/kg-d)	1/(mg/kg-d)		mg/kg-d			mg/kg-d			mg/kg-d			mg/kg-d			
1,2,4-TRIMETHYLBENZENE	2.87E-01	NA	NA	NC	NA	NA	1	NA	NA	NC	NA	NA	1	NA	NA	NA	NA
1,2-DICHLOROETHANE	2.06E-02	0.02	0.091	1	3.53E-10	3.22E-11	1	3.53E-09	1.77E-07	0.1	4.21E-10	3.84E-11	0.1	4.21E-09	2.11E-07	7.05E-11	3.87E-07
1,3,5-TRIMETHYLBENZENE	1.15E-01	0.01	NA	NC	NA	NA	1	1.97E-08	1.97E-06	NC	NA	NA	1	2.34E-07	2.34E-05	NA	2.54E-05
4-CHLOROANILINE	6.36E-01	0.004	0.2	1	1.09E-08	2.18E-09	1	1.09E-07	2.73E-05	0.08	1.04E-08	2.08E-09	0.1	1.30E-07	3.25E-05	4.26E-09	5.98E-05
4-ISOPROPYLTOLUENE	2.59E-02	NA	NA	NC	NA	NA	1	NA	NA	NC	NA	NA	0.12	NA	NA	NA	NA
ACETONE	2.05E+00	0.9	NA	NC	NA	NA	1	3.52E-07	3.91E-07	NC	NA	NA	0.1	4.20E-07	4.67E-07	NA	8.58E-07
ANTIMONY	2.18E+00	0.0004	NA	NC	NA	NA	1	3.74E-07	9.35E-04	NC	NA	NA	0.1	4.46E-07	1.12E-03	NA	2.05E-03
AROCLOR-1260	1.33E-01	0.00002	2	0.85	1.94E-09	3.88E-09	0.85	1.94E-08	9.70E-04	0.16	4.36E-09	8.71E-09	0.16	4.36E-08	2.18E-03	1.26E-08	3.15E-03
ARSENIC	4.73E+00	0.0003	1.5	1	8.13E-08	1.22E-07	1	8.13E-07	2.71E-03	0.03	2.91E-08	4.36E-08	0.03	2.91E-07	9.69E-04	1.66E-07	3.68E-03
BARIUM	2.49E+02	0.2	NA	NC	NA	NA	1	4.27E-05	2.13E-04	NC	NA	NA	0.05	2.54E-05	1.27E-04	NA	3.41E-04
BENZO(B)FLUORANTHENE	2.87E-01	0.03	0.73	0.28	1.38E-09	1.01E-09	0.28	1.38E-08	4.60E-07	0.02	1.18E-09	8.59E-10	0.02	1.18E-08	3.92E-07	1.87E-09	8.52E-07
BENZO(G,H,I)PERYLENE	2.01E-01	0.03	NA	NC	NA	NA	0.36	1.25E-08	4.15E-07	NC	NA	NA	0.1	4.13E-08	1.38E-06	NA	1.79E-06
BENZO(K)FLUORANTHENE	3.00E-01	0.03	0.073	0.28	1.44E-09	1.05E-10	0.28	1.44E-08	4.81E-07	0.02	1.23E-09	8.98E-11	0.02	1.23E-08	4.10E-07	1.95E-10	8.91E-07
BERYLLÌÚM	5.43E-01	0.002	NA	NC	NA	NA	1	9.32E-08	4.66E-05	0.03	NA	NA	0.03	3.34E-08	1.67E-05	NA	6.33E-05
C11-C22 AROMATICS	2.78E+01	0.03	NA	NC	NA	NA	0.36	1.72E-06	5.73E-05	NC	NA	NA	0.1	5.70E-06	1.90E-04	NA	2.47E-04
C19-C36 ALIPHATICS	4.81E+01	2	NA	NC	NA	NA	1	8.25E-06	4.13E-06	NC	NA	NA	0.1	9.84E-06	4.92E-06	NA	9.05E-06
C5-C8 ALIPHATICS	2.04E+00	0.04	NA	NC	NA	NA	1	3.50E-07	8.76E-06	NC	NA	NA	1	4.18E-06	1.05E-04	NA	1.13E-04
C9-C10 AROMATICS	5.17E+00	0.03	NA	NC	NA	NA	1	8.88E-07	2.96E-05	NC	NA	NA	0.5	5.30E-06	1.77E-04	NA	2.06E-04
C9-C18 ALIPHATICS	4.04E+00	0.1	NA	NC	NA	NA	1	6.94E-07	6.94E-06	NC	NA	NA	0.5	4.14E-06	4.14E-05	NA	4.83E-05
CADMIUM	9.26E-01	0.0005	NA	NC	NA	NA	1_	1.59E-07	3.18E-04	NC	NA	NA	0.14	2.66E-07	5.31E-04	NA	8.49E-04

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Attachment D
Exposure and Risk Estimates Associated With Soil Contact
Bird Machine Company - DDA
100 Neponset St, Walpole MA
Trespasser Older Child
Surface Soil
DDA current
0-3 feet

Receptor:	Trespasser Older Child	▼
Medium:	Surface Soil	•
Exposure Area:	DDA current	▼
Depth:	0-3 feet	<u></u>
Duration:	Chronic	▼

Parameter	Definition	Units	Value	Comment	
IRsoil	Soil Ingestion Rate	mg/d	50		
SA	Soil Dermal Contact Skin Exposed	cm2/d	4260		
AF	Soil Dermal Contact Adherence Rate	mg/cm2	0.14		
EF	Soil Exposure Frequency	d/y	50		
EP	Soil Exposure Period - Cancer	У	7		
EP	Soil Exposure Period - Non-Cancer	y	7		
ATc	Soil Averaging Time - Cancer	d	25550		
ATn	Soil Averaging Time - Non-Cancer	d	2555		
BW	Body Weight	kg	39.9		
CF	Conversion Factor	kg/mg	0.000001		

$ADD_{ing} = \frac{C_{soil} \times CF \times IR_{soil} \times RAF_{os} \times EF \times EP}{AT \times BW}$
$ ADD _{der} = \frac{C_{soil} \times CF \times SA \times AF \times RAF _{ds} \times EF \times EP}{AT - BW}$
$AT \times BW$
$HI_{ing} = \frac{ADD_{ing}}{RfD}$
$HI_{der} = \frac{ADD_{der}}{RfD}$
$HI = HI_{ing} + HI_{der}$
$ Risk_{ing} = ADD_{ing} \times CSF$
$ Risk _{der} = ADD _{der} \times CSF$
$Risk = Risk_{ing} + Risk_{der}$

	EPC			Incidental In	gestion					Dermal Co	ontact					Total	•
	Surface Soil																
	DDA current																
	0-3 feet	RfD	CSF	RAFosc	ADDing-c	Risking	RAFosnc	ADDing-nc	Hling	RAFdsc	ADDder-c	Riskder	RAFdsnc	ADDder-nc	Hlder	Risk (Soil)	HI (Soil)
Compound	(mg/kg)	(mg/kg-d)	1/(mg/kg-d)		mg/kg-d			mg/kg-d			mg/kg-d			mg/kg-d			
CHROMIUM	1.89E+02	1.5	NA	NC	NA	NA	1	3.25E-05	2.17E-05	NC	NA	NA	0.04	1.55E-05	1.03E-05	NA	3.20E-05
COPPER	5.60E+01	0.04	NA NA	NC	NA	NA	1	9.61E-06	2.17E-03 2.40E-04	NC	NA NA	NA	0.04	1.15E-04	2.87E-03	NA NA	3.20L-03 3.11E-03
DIBENZ(A,H)ANTHRACENE	2.01E-01	0.04	7.3	0.28	9.68E-10	7.07E-09	0.28	9.68E-09	3.23E-07	0.02	8.25E-10	6.02E-09	0.02	8.25E-09	2.07E-03 2.75E-07	1.31E-08	5.11L-03 5.98E-07
ETHYLBENZENE			7.3 0.11	0.20			0.20										
FLUORANTHENE	3.33E-02	0.1		NC.	5.71E-10	6.29E-11	0.20	5.71E-09	5.71E-08	0.08 NC	5.45E-10	6.00E-11	0.2	1.36E-08	1.36E-07	1.23E-10	1.93E-07
	6.30E-01	0.04	NA	NC	NA	NA	0.36	3.89E-08	9.73E-07	_	NA	NA	0.1	1.29E-07	3.22E-06	NA	4.20E-06
LEAD	1.01E+02	0.00075	NA	NC	NA	NA	0.5	8.64E-06	1.15E-02	NC	NA	NA	0.006	1.24E-06	1.65E-03	NA	1.32E-02
M,P-XYLENES	1.78E-01	0.2	NA	NC	NA	NA	1	3.06E-08	1.53E-07	NC	NA	NA	0.12	4.38E-08	2.19E-07	NA	3.72E-07
MERCURY	4.19E-01	0.0003	NA	NC	NA	NA	1	7.19E-08	2.40E-04	NC	NA	NA	0.05	4.29E-08	1.43E-04	NA	3.83E-04
N-BUTYLBENZENE	2.27E-02	0.05	NA	NC	NA	NA	1	3.90E-09	7.79E-08	NC	NA	NA	1	4.65E-08	9.29E-07	NA	1.01E-06
NICKEL	2.25E+02	0.02	NA	NC	NA	NA	1	3.86E-05	1.93E-03	NC	NA	NA	0.35	1.61E-04	8.06E-03	NA	9.99E-03
N-PROPYLBENZENE	4.11E-02	0.1	NA	NC	NA	NA	1	7.06E-09	7.06E-08	NC	NA	NA	1	8.42E-08	8.42E-07	NA	9.13E-07
O-XYLENE	7.04E-02	0.2	NA	NC	NA	NA	1	1.21E-08	6.04E-08	NC	NA	NA	0.12	1.73E-08	8.65E-08	NA	1.47E-07
PHENANTHRENE	5.35E-01	0.03	NA	NC	NA	NA	0.36	3.31E-08	1.10E-06	NC	NA	NA	0.1	1.10E-07	3.65E-06	NA	4.75E-06
PYRENE	6.02E-01	0.03	NA	NC	NA	NA	0.36	3.72E-08	1.24E-06	NC	NA	NA	0.1	1.23E-07	4.11E-06	NA	5.35E-06
SILVER	1.16E+00	0.005	NA	NC	NA	NA	1	2.00E-07	3.99E-05	NC	NA	NA	0.25	5.96E-07	1.19E-04	NA	1.59E-04
TERT-BUTYLBENZENE	3.54E-02	NA	NA	NC	NA	NA	1	NA	NA	NC	NA	NA	1	NA	NA	NA	NA
TOLUENE	4.00E-02	0.08	NA	NC	NA	NA	1	6.87E-09	8.59E-08	NC	NA	NA	0.12	9.83E-09	1.23E-07	NA	2.09E-07
VANADIUM	3.83E+01	0.009	NA	NC	NA	NA	1	6.57E-06	7.30E-04	NC	NA	NA	0.03	2.35E-06	2.61E-04	NA	9.91E-04
ZINC	1.69E+02	0.3	NA	NC	NA	NA	1	2.90E-05	9.68E-05	NC	NA	NA	0.02	6.93E-06	2.31E-05	NA	1.20E-04
2,3,7,8-TCDD TEQ	2.75E-05	0.000000001	130000	1	4.73E-13	6.15E-08	1	4.73E-12	4.73E-03	0.2	1.13E-12	1.47E-07	1	5.64E-11	5.64E-02	2.08E-07	6.11E-02
						1.98E-07			2.49E-02			2.08E-07			7.51E-02	4.06E-07	9.99E-02
						1.300-07			2. 4 3E-02			2.00E-07			1.01E-02	4.00E-07	∌.∌∌E-UZ

NA - Not available NC - Not calculated ND - Not detected

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Attachment D
Exposure and Risk Estimates Associated With Inhalation of Soil-Derived Particulates in Air
Bird Machine Company - DDA
100 Neponset St, Walpole MA
Trespasser Older Child
Ambient Air
DDA current
0-3 feet

Receptor:	Trespasser Older Child	•
Medium of Origin:	Surface Soil	•
Exposure Medium:	Ambient Air	▼
Exposure Area:	DDA current	_
Depth:	0-3 feet	•
Duration:	Chronic	_

Parameter	Definition	Units	Value	Comment	
IRair	Outdoor Air Inhalation Rate - Particulates	m3/hr	0.94		
PEF	Outdoor Air PM10 - Particulates	ug/m3	32		
ET	Outdoor Air Exposure Time - Particulates	hr/d	2		
EF	Outdoor Air Exposure Frequency - Particulates	d/y	50		
EP	Outdoor Air Exposure Period - Cancer - Particu	ıle y	7		
EP	Outdoor Air Exposure Period - Non-Cancer - P	ar y	7		
ATc	Outdoor Air Averaging Time - Cancer - Particulates	d	25550		
ATn	Outdoor Air Averaging Time - Non-Cancer - Particulates	d	2555		
BW	Body Weight	kg	39.9		
С	Conversion Factor	ug/mg	1000		

$$C_{air} = C_{soil} \times PM_{10} \times 1 \times 10^{-9} \, kg / ug$$

$$ADD_{inh-gi} = \frac{C_{air} \times 1.5 \times IR_{air} \times RAF_{i} \times ET \times EF \times EP}{AT \times BW}$$

$$ADD_{inh} = \frac{C_{air} \times 0.5 \times IR_{air} \times RAF_{i} \times ET \times EF \times EP}{AT \times BW}$$

$$ADE_{inh} = \frac{ADD_{inh} \times 70 \, kg}{20 \, m^{3} / d}$$

$$HI = \frac{ADD_{inh-gi}}{RfD} + \frac{ADE_{inh}}{RfC}$$

$$Risk = (ADD_{inh-gi} \times CSF) + (ADE_{inh} \times URF \times C)$$

	EPC	EPC																		
	Surface Soil																		Risk	HI
	DDA current																		(Particulates	(Particulate
	0-3 feet	Fugitive Dust	RfC	URF	RfD	CSF	RAFic	ADD-inhc	ADE-c	Riskinh	ADD-ingc	Risking	RAFinc	ADD-inhnc	ADE-nc	Hlinh	ADD-ingnc	Hling	in Air)	in Air)
Compound	(mg/kg)	(mg/m3)	(mg/m3)	1/(ug/m3)	(mg/kg-d)	1/(mg/kg-d)		mg/kg-d	mg/m3					mg/kg-d	mg/m3					
1,2,4-TRIMETHYLBENZENE	2.87E-01	9.18E-09	0.007	NA	NA	NA	NC	NA	NA	NA	NA	NA	1	2.96E-11	1.04E-10	1.48E-08	NA	NA	NA	1.48E-08
1,2-DICHLOROETHANE	2.06E-02	6.59E-10	0.055	0.000026	0.02	0.091	1	2.126E-13	7.439E-13	1.93E-14	6.377E-13	5.81E-14	1	2.13E-12	7.44E-12	1.35E-10	6.38E-12	3.19E-10	7.74E-14	4.54E-10
1,3,5-TRIMETHYLBENZENE	1.15E-01	3.66E-09	NA	NA	0.01	NA	NC	NA	NA	NA	NA	NA	1	NA	NA	NA	3.55E-11	3.55E-09	NA	3.55E-09
4-CHLOROANILINE	6.36E-01	2.03E-08	0.014	NA	0.004	0.2	1	NA	NA	NA	1.969E-11	3.94E-12	1	6.56E-11	2.30E-10	1.64E-08	1.97E-10	4.92E-08	3.94E-12	6.56E-08
4-ISOPROPYLTOLUENE	2.59E-02	8.28E-10	NA	NA	NA	NA	NC	NA	NA	NA	NA	NA	1	NA	NA	NA	NA	NA	NA	NA
ACETONE	2.05E+00	6.56E-08	0.8	NA	0.9	NA	NC	NA	NA	NA	NA	NA	1	2.12E-10	7.41E-10	9.27E-10	6.35E-10	7.06E-10	NA	1.63E-09
ANTIMONY	2.18E+00	6.97E-08	0.01	NA	0.0004	NA	NC	NA	NA	NA	NA	NA	1	2.25E-10	7.87E-10	7.87E-08	6.75E-10	1.69E-06	NA	1.77E-06
AROCLOR-1260	1.33E-01	4.25E-09	0.00002	0.0001	0.00002	2	1	1.373E-12	4.805E-12	4.81E-13	4.119E-12	8.24E-12	1	1.37E-11	4.81E-11	2.40E-06	4.12E-11	2.06E-06	8.72E-12	4.46E-06
ARSENIC	4.73E+00	1.51E-07	0.0000025	0.0043	0.0003	1.5	1	4.889E-11	1.711E-10	7.36E-10	1.467E-10	2.20E-10	1	4.89E-10	1.71E-09	6.84E-04	1.47E-09	4.89E-06	9.56E-10	6.89E-04
BARIUM	2.49E+02	7.95E-06	0.0005	NA	0.2	NA	NC	NA	NA	NA	NA	NA	1	2.57E-08	8.98E-08	1.80E-04	7.70E-08	3.85E-07	NA	1.80E-04
BENZO(B)FLUORANTHENE	2.87E-01	9.19E-09	0.05	0.00011	0.03	0.73	1	2.966E-12	1.038E-11	1.14E-12	8.897E-12	6.49E-12	1	2.97E-11	1.04E-10	2.08E-09	8.90E-11	2.97E-09	7.64E-12	5.04E-09
BENZO(G,H,I)PERYLENE	2.01E-01	6.45E-09	0.05	NA	0.03	NA	NC	NA	NA	NA	NA	NA	1	2.08E-11	7.28E-11	1.46E-09	6.24E-11	2.08E-09	NA	3.54E-09
BENZO(K)FLUORANTHENE	3.00E-01	9.61E-09	0.05	0.00011	0.03	0.073	1	3.101E-12	1.085E-11	1.19E-12	9.304E-12	6.79E-13	1	3.10E-11	1.09E-10	2.17E-09	9.30E-11	3.10E-09	1.87E-12	5.27E-09
BERYLLIÚM	5.43E-01	1.74E-08	0.00002	0.0024	0.002	NA	1	5.607E-12	1.962E-11	4.71E-11	NA	NA	1	5.61E-11	1.96E-10	9.81E-06	1.68E-10	8.41E-08	4.71E-11	9.90E-06
C11-C22 AROMATICS	2.78E+01	8.90E-07	0.05	NA	0.03	NA	NC	NA	NA	NA	NA	NA	1	2.87E-09	1.01E-08	2.01E-07	8.62E-09	2.87E-07	NA	4.88E-07
C19-C36 ALIPHATICS	4.81E+01	1.54E-06	NA	NA	2	NA	NC	NA	NA	NA	NA	NA	1	NA	NA	NA	1.49E-08	7.45E-09	NA	7.45E-09
C5-C8 ALIPHATICS	2.04E+00	6.53E-08	0.2	NA	0.04	NA	NC	NA	NA	NA	NA	NA	1	2.11E-10	7.38E-10	3.69E-09	6.33E-10	1.58E-08	NA	1.95E-08
C9-C10 AROMATICS	5.17E+00	1.66E-07	0.05	NA	0.03	NA	NC	NA	NA	NA	NA	NA	1	5.34E-10	1.87E-09	3.74E-08	1.60E-09	5.34E-08	NA	9.09E-08
C9-C18 ALIPHATICS	4.04E+00	1.29E-07	0.2	NA	0.1	NA	NC	NA	NA	NA	NA	NA	1	4.17E-10	1.46E-09	7.30E-09	1.25E-09	1.25E-08	NA	1.98E-08
CADMIUM	9.26E-01	2.96E-08	0.00002	0.0018	0.0005	NA	1	9.568E-12	3.349E-11	6.03E-11	NA	NA	1	9.57E-11	3.35E-10	1.67E-05	2.87E-10	5.74E-07	6.03E-11	1.73E-0
CHROMIUM	1.89E+02	6.06E-06	0.0001	NA	1.5	NA	NC	NA	NA	NA	NA	NA	1	1.96E-08	6.85E-08	6.85E-04	5.87E-08	3.91E-08	NA	6.85E-04
COPPER	5.60E+01	1.79E-06	NA	NA	0.04	NA	1	NA	NA	NA	NA	NA	1	NA	NA	NA	1.73E-08	4.34E-07	NA	4.34E-07
DIBENZ(A.H)ANTHRACENE	2.01E-01	6.45E-09	0.05	0.0012	0.03	7.3	1	2.081E-12	7 282F-12	8 74F-12	6.242E-12	4.56E-11	1	2.08E-11	7.28E-11	1.46E-09	6.24E-11	2.08E-09	5.43E-11	3.54E-09
ETHYLBENZENE	3.33E-02	1.07E-09	1	0.0000025	0.1	0.11	1	3.437E-13	1.203E-12	3.00E-15	1.031E-12	1.13E-13	l i	3.44E-12	1.20E-11	1.20E-11	1.03E-11	1.03E-10	1.16E-13	1.15E-10

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Attachment D
Exposure and Risk Estimates Associated With Inhalation of Soil-Derived Particulates in Air
Bird Machine Company - DDA
100 Neponset St, Walpole MA
Trespasser Older Child
Ambient Air
DDA current
0-3 feet

Receptor:	Trespasser Older Child	
Medium of Origin:	Surface Soil	•
Exposure Medium:	Ambient Air	
Exposure Area:	DDA current	▼
Depth:	0-3 feet	•
Duration:	Chronic	▼

Parameter	Definition	Units	Value	Comment	
IRair	Outdoor Air Inhalation Rate - Particulates	m3/hr	0.94		
PEF	Outdoor Air PM10 - Particulates	ug/m3	32		
ET	Outdoor Air Exposure Time - Particulates	hr/d	2		
EF	Outdoor Air Exposure Frequency - Particulates	d/y	50		
EP	Outdoor Air Exposure Period - Cancer - Particu	la y	7		
EP	Outdoor Air Exposure Period - Non-Cancer - Pa	ar y	7		
ATc	Outdoor Air Averaging Time - Cancer - Particulates	d	25550		
ATn	Outdoor Air Averaging Time - Non-Cancer - Particulates	d	2555		
BW	Body Weight	kg	39.9		
С	Conversion Factor	ug/mg	1000		

$$C_{air} = C_{soil} \times PM_{10} \times 1 \times 10^{-9} \, kg / ug$$

$$ADD_{inh-gi} = \frac{C_{air} \times 1.5 \times IR_{air} \times RAF_{i} \times ET \times EF \times EP}{AT \times BW}$$

$$ADD_{inh} = \frac{C_{air} \times 0.5 \times IR_{air} \times RAF_{i} \times ET \times EF \times EP}{AT \times BW}$$

$$ADE_{inh} = \frac{ADD_{inh} \times 70 \, kg}{20 \, m^{3} / d}$$

$$HI = \frac{ADD_{inh-gi}}{RfD} + \frac{ADE_{inh}}{RfC}$$

$$Risk = (ADD_{inh-gi} \times CSF) + (ADE_{inh} \times URF \times C)$$

	EPC	EPC																	Dist	
	Surface Soil DDA current																		Risk (Particulates	HI (Particulates
	0-3 feet	Fugitive Dust	RfC	URF	RfD	CSF	RAFic	ADD-inhc	ADE-c	Riskinh	ADD-ingc	Risking	RAFinc	ADD-inhnc	ADE-nc	Hlinh	ADD-ingnc	Hlina	in Air)	in Air)
Compound	(mg/kg)	(mg/m3)	(mg/m3)	1/(ug/m3)		1/(mg/kg-d)	101110	mg/kg-d	mg/m3		, as a migo	ruotung	70 1110	mg/kg-d	mg/m3		7.22 mgno		,	,
FLUORANTHENE	6.30E-01	2.01E-08	0.05	NA	0.04	NA	NC	NA	NA	NA	NA	NA	1	6.50E-11	2.28E-10	4.55E-09	1.95E-10	4.88E-09	NA	9.43E-09
LEAD	1.01E+02	3.22E-06	0.001	NA	0.00075	NA	NC	NA	NA	NA	NA	NA	1	1.04E-08	3.64E-08	3.64E-05	3.12E-08	4.16E-05	NA	7.80E-05
M,P-XYLENES	1.78E-01	5.70E-09	0.1	NA	0.2	NA	NC	NA	NA	NA	NA	NA	1	1.84E-11	6.44E-11	6.44E-10	5.52E-11	2.76E-10	NA	9.20E-10
MERCURY	4.19E-01	1.34E-08	0.0003	NA	0.0003	NA	NC	NA	NA	NA	NA	NA	1	4.33E-11	1.51E-10	5.05E-07	1.30E-10	4.33E-07	NA	9.38E-07
N-BUTYLBENZENE	2.27E-02	7.26E-10	NA	NA	0.05	NA	NC	NA	NA	NA	NA	NA	1	NA	NA	NA	7.03E-12	1.41E-10	NA	1.41E-10
NICKEL	2.25E+02	7.20E-06	0.001	0.00026	0.02	NA	1	2.323E-09	8.129E-09	2.11E-09	NA	NA	1	2.32E-08	8.13E-08	8.13E-05	6.97E-08	3.48E-06	2.11E-09	8.48E-05
N-PROPYLBENZENE	4.11E-02	1.32E-09	1	NA	0.1	NA	NC	NA	NA	NA	NA	NA	1	4.25E-12	1.49E-11	1.49E-11	1.27E-11	1.27E-10	NA	1.42E-10
O-XYLENE	7.04E-02	2.25E-09	0.1	NA	0.2	NA	NC	NA	NA	NA	NA	NA	1	7.27E-12	2.55E-11	2.55E-10	2.18E-11	1.09E-10	NA	3.64E-10
PHENANTHRENE	5.35E-01	1.71E-08	0.05	NA	0.03	NA	NC	NA	NA	NA	NA	NA	1	5.53E-11	1.93E-10	3.87E-09	1.66E-10	5.53E-09	NA	9.39E-09
PYRENE	6.02E-01	1.93E-08	0.05	NA	0.03	NA	NC	NA	NA	NA	NA	NA	1	6.22E-11	2.18E-10	4.35E-09	1.86E-10	6.22E-09	NA	1.06E-08
SILVER	1.16E+00	3.72E-08	0.00014	NA	0.005	NA	NC	NA	NA	NA	NA	NA	1	1.20E-10	4.21E-10	3.00E-06	3.60E-10	7.21E-08	NA	3.08E-06
TERT-BUTYLBENZENE	3.54E-02	1.13E-09	NA	NA	NA	NA	NC	NA	NA	NA	NA	NA	1	NA	NA	NA	NA	NA	NA	NA
TOLUENE	4.00E-02	1.28E-09	5	NA	0.08	NA	NC	NA	NA	NA	NA	NA	1	4.13E-12	1.45E-11	2.89E-12	1.24E-11	1.55E-10	NA	1.58E-10
VANADIUM	3.83E+01	1.22E-06	0.001	NA	0.009	NA	NC	NA	NA	NA	NA	NA	1	3.95E-09	1.38E-08	1.38E-05	1.19E-08	1.32E-06	NA	1.52E-05
ZINC	1.69E+02	5.41E-06	0.0014	NA	0.3	NA	NC	NA	NA	NA	NA	NA	1	1.75E-08	6.12E-08	4.37E-05	5.24E-08	1.75E-07	NA	4.39E-05
2,3,7,8-TCDD TEQ	2.75E-05	8.81E-13	0.00000004	38	1E-09	130000	1	2.845E-16	9.956E-16	3.78E-11	8.534E-16	1.11E-10	1	2.84E-15	9.96E-15	2.49E-07	8.53E-15	8.53E-06	1.49E-10	8.78E-06
Total										3.01E-09		3.96E-10				1.76E-03		6.62E-05	3.40E-09	1.82E-03

NA - Not available NC - Not calculated ND - Not detected

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Attachment D Human Health Risk Calculations for Soil - Asbestos Bird Machine Company - Walpole, MA Human Health and Environmental Risk Characterization

Calculations: OHMsoil fibers/g	PM10 mg/m3	C g-m3/mg-ml	OHMair fibers/ml	
8.29E+07	0.032	1.00E-09	2.65E-03	
OHMair fibers/ml	EF days/year	EP years	AP days	ADEair fibers/ml
2.65E-03	50	7	25550	3.63E-05
ADEair fibers/ml	IUR ml/fiber	ELCR		
3.63E-05	0.23	8.4E-06		

Equations:

$$OHM_{air} = OHM_{soil} \times PM_{10} \times C$$

$$ADC_{air} = \frac{OHM_{air} \times EF \times EP}{AP}$$

ELCR = ADEair x IUR

Notes:

OHM - Oil and/or hazardous materials

fibers/g - fibers per gram

PM10 - particulate matter measuring 10 micrometers or less

ml - milliliter

C - conversion factor

EF - exposure frequency

EP - exposure period

AP - averging period

ADE - average daily exposure

IUR - inhalation unit risk

ELCR - excess lifetime cancer risk

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DD-MW-002

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 8E-07 HI (all chemicals) = 7E-02

6 4 : 466	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQ_{ing}	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	2.6E+00	6.7E-07	1.3E-07	7.6E-10	8.0E-07	7.5E-03	1.5E-03	2.2E-04	9.2E-03
Barium	1.9E+01					5.6E-03	3.3E-05		5.6E-03
Chromium (total)	1.9E+00					3.7E-02	3.7E-03		4.1E-02
Nickel	1.2E+00					3.5E-03	3.8E-05		3.6E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	5.0E-01					3.3E-03	3.6E-04		3.6E-03
Zinc	1.2E+00					2.4E-04	1.7E-06		2.4E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-201

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 5E-07 HI (all chemicals) = 2E-01

Contaminant of Concern	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQing	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	1.5E+00	4.0E-07	7.9E-08	7.6E-10	4.8E-07	4.5E-03	8.9E-04	2.2E-04	5.6E-03
Barium	4.6E+01					1.3E-02	3.3E-05		1.3E-02
Chromium (total)	2.8E+00					5.5E-02	3.7E-03		5.9E-02
Nickel	1.3E+00					3.7E-03	3.8E-05		3.7E-03
Selenium	7.3E+00					8.6E-02	5.3E-05		8.6E-02
Vanadium	3.4E-01					2.2E-03	3.6E-04		2.6E-03
Zinc	2.5E+00	·	·			4.9E-04	1.7E-06		4.9E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-203

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 1E-07 HI (all chemicals) = 3E-02

Contaminant of Consum	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQ _{ing}	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	4.3E-01	1.1E-07	2.2E-08	7.6E-10	1.4E-07	1.3E-03	2.5E-04	2.2E-04	1.7E-03
Barium	1.1E+01					3.2E-03	3.3E-05		3.3E-03
Chromium (total)	6.3E-01					1.2E-02	3.7E-03		1.6E-02
Nickel	1.2E+00					3.5E-03	3.8E-05		3.6E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	2.5E-01					1.6E-03	3.6E-04		2.0E-03
Zinc	5.3E+00					1.0E-03	1.7E-06		1.0E-03

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-204

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 1E-07 HI (all chemicals) = 4E-02

Contonin ant of Consum	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQ _{ing}	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	4.5E-01	1.2E-07	2.4E-08	7.6E-10	1.4E-07	1.3E-03	2.6E-04	2.2E-04	1.8E-03
Barium	2.0E+01					5.9E-03	3.3E-05		5.9E-03
Chromium (total)	5.0E-01					9.8E-03	3.7E-03		1.3E-02
Nickel	1.2E+00					3.5E-03	3.8E-05		3.6E-03
Selenium	1.0E+00					1.2E-02	5.3E-05		1.2E-02
Vanadium	5.0E-01					3.3E-03	3.6E-04		3.6E-03
Zinc	2.5E+00					4.9E-04	1.7E-06		4.9E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-205

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 8E-07 HI (all chemicals) = 4E-02

Contaminant of Concern	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQ _{ing}	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	2.6E+00	6.7E-07	1.3E-07	7.6E-10	8.0E-07	7.5E-03	1.5E-03	2.2E-04	9.2E-03
Barium	7.6E+00					2.2E-03	3.3E-05		2.3E-03
Chromium (total)	5.2E-01					1.0E-02	3.7E-03		1.4E-02
Nickel	5.9E-01					1.7E-03	3.8E-05		1.8E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	2.9E-01					1.9E-03	3.6E-04		2.3E-03
Zinc	1.6E+00					3.1E-04	1.7E-06		3.2E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-206

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 2E-07 HI (all chemicals) = 3E-02

Contaminant of Concern	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQ _{ing}	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	4.9E-01	1.3E-07	2.6E-08	7.6E-10	1.5E-07	1.4E-03	2.9E-04	2.2E-04	2.0E-03
Barium	1.6E+01					4.7E-03	3.3E-05		4.7E-03
Chromium (total)	5.1E-01					1.0E-02	3.7E-03		1.4E-02
Nickel	6.9E-01					2.0E-03	3.8E-05		2.1E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	5.0E-01					3.3E-03	3.6E-04		3.6E-03
Zinc	1.9E+00					3.7E-04	1.7E-06		3.7E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-207

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 2E-07 HI (all chemicals) = 5E-02

Contaminant of Concern	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQing	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	5.1E-01	1.3E-07	2.7E-08	7.6E-10	1.6E-07	1.5E-03	3.0E-04	2.2E-04	2.0E-03
Barium	3.1E+01					9.1E-03	3.3E-05		9.2E-03
Chromium (total)	1.3E+00					2.5E-02	3.7E-03		2.9E-02
Nickel	4.3E-01					1.3E-03	3.8E-05		1.3E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	5.0E-01					3.3E-03	3.6E-04		3.6E-03
Zinc	1.9E+00	·	·			3.7E-04	1.7E-06		3.7E-04

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient

DD-MW-208

Resident - Drinking Water: Table RW-1

Exposure Point Concentration (EPC) and Risk

Based on Resident Ages 1-31 (Cancer) and 1-8 (Noncancer)

ShortForm Version 4-06 Vlookup Version v0808

ELCR (all chemicals) = 5E-07 HI (all chemicals) = 6E-02

Contaminant of Concern	EPC	ELCR	ELCR	ELCR			Chronic		
Contaminant of Concern	(ug/L)	ingestion	dermal	inhalation	ELCR _{total}	HQing	HQ _{derm}	HQ _{inh}	HQ _{total}
Bis(2-ethylhexyl)phthalate	1.6E+00	4.2E-07	8.4E-08	7.6E-10	5.0E-07	4.7E-03	9.4E-04	2.2E-04	5.9E-03
Barium	3.1E+01					9.1E-03	3.3E-05		9.2E-03
Chromium (total)	3.6E-01					7.1E-03	3.7E-03		1.1E-02
Nickel	2.0E+00					5.9E-03	3.8E-05		5.9E-03
Selenium	5.0E-01					5.9E-03	5.3E-05		5.9E-03
Vanadium	5.2E-01					3.4E-03	3.6E-04		3.8E-03
Zinc	1.1E+02	·	·			2.2E-02	1.7E-06		2.2E-02

Notes:

EPC - The concentration detected in the monitoring well during the 2008 sampling round. One-half of the reporting limit was used if the analyte was not detected.

ug/L - micrograms per liter

ELCR - excess lifetime cancer risk

HQ - hazard quotient



Attachment F-1 Table 1A - Evaluation of Potential Risk to Red Fox Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterizatrion

Site: DDA Red Fox Receptor: Pathway: Soil ingestion Consumption of small mammals

Parameter	Value	
Small Mammal Ingestion Rate (kg ww/d)	0.510	Assume 100% of diet.
Total Dietary Intake (kg ww/d)	0.510	USEPA (1993).
Soil Ingestion Rate (kg dw/day)	0.01122	Calc. from Beyer (1994)
Body Weight (kg)	4.54	USEPA (1993). Mean of male and female BWs in spring and fall.
Sm. Mammal Dry wt./wet wt. CF	0.3	USEPA (1999)
Home range (ha)	1038	USEPA (1993)
Area Use Factor	0.00178	1.85. ha exposure area

	Surface Soil											
			Conc. in		Intake from							
NOAEL-			small	Intake	small							
based TRV	Conc. in soil		mammals	from soil	mammals		TQ-small					
(mg/kg-d)	(mg/kg)	BTF	(mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	mammals	Total TQ				
3.79E+00	2.24E+00	(a)	3.5E-02	9.86E-06	2.09E-06	2.60E-06	5.52E-07	3.2E-06				
7.04E-01	4.84E+00	(a)	6.9E-02	2.13E-05	4.13E-06	3.03E-05	5.87E-06	3.6E-05				
1.48E+01	2.62E+02	(a)	4.8E-01	1.16E-03	2.88E-05	7.83E-05	1.95E-06	8.0E-05				
5.29E-01	9.76E-01	(a)	2.8E-01	4.30E-06	1.69E-05	8.12E-06	3.19E-05	4.0E-05				
6.85E-01	2.00E+02	(a)	1.7E+01	8.82E-04	1.01E-03	1.29E-03	1.47E-03	2.8E-03				
7.12E+00	5.60E+01	(a)	2.4E+00	2.47E-04	1.44E-04	3.47E-05	2.03E-05	5.5E-05				
1.16E+01	1.06E+02	(a)	8.1E+00	4.69E-04	4.89E-04	4.04E-05	4.21E-05	8.3E-05				
6.85E-01	4.43E-01	7.5E-06	3.3E-06	1.95E-06	6.67E-10	2.85E-06	9.74E-10	2.9E-06				
2.19E+00	2.40E+02	(a)	1.0E+01	1.06E-03	6.03E-04	4.81E-04	2.75E-04	7.6E-04				
1.69E+00	3.89E+01	1.2E-02	4.8E-01	1.71E-04	2.87E-05	1.01E-04	1.70E-05	1.2E-04				
2.15E+01	1.78E+02	(a)	1.1E+02	7.83E-04	6.80E-03	3.64E-05	3.16E-04	3.5E-04				
1.60E-02	1.38E-01	5.8E-05	8.0E-06	6.07E-07	1.61E-09	3.79E-05	1.00E-07	3.8E-05				
5.27E-07	2.76E-05	7.8E-05	2.2E-09	1.22E-10	4.31E-13	2.31E-04	8.19E-07	2.3E-04				
5.12E+00	2.99E-01	0.0E+00	0.0E+00	1.32E-06	0.00E+00	2.57E-07	0.00E+00	2.6E-07				
5.12E+00	2.08E-01	0.0E+00	0.0E+00	9.17E-07	0.00E+00	1.79E-07	0.00E+00	1.8E-07				
5.12E+00	3.13E-01	0.0E+00	0.0E+00	1.38E-06	0.00E+00	2.69E-07	0.00E+00	2.7E-07				
5.12E+00	2.08E-01	0.0E+00	0.0E+00	9.17E-07	0.00E+00	1.79E-07	0.00E+00	1.8E-07				
5.12E+00	6.34E-01	0.0E+00	0.0E+00	2.79E-06	0.00E+00	5.45E-07	0.00E+00	5.4E-07				
2.14E+01	2.04E+00	1.0E+00	2.0E+00	9.00E-06	1.23E-04	4.19E-07	5.72E-06	6.1E-06				
2.14E+01	5.17E+00	1.0E+00	5.2E+00	2.28E-05	3.11E-04	1.07E-06	1.45E-05	1.6E-05				
5.27E+01	1.41E+00	1.0E+00	1.4E+00	6.21E-06	8.46E-05	1.18E-07	1.61E-06	1.7E-06				
2.14E+01	2.97E+01	1.0E+00	3.0E+01	1.31E-04	1.78E-03	6.11E-06	8.33E-05	8.9E-05				
1.05E+02	5.14E+01	1.0E+00	5.1E+01	2.26E-04	3.09E-03	2.15E-06	2.93E-05	3.1E-05				
	based TRV (mg/kg-d) 3.79E+00 7.04E-01 1.48E+01 5.29E-01 6.85E-01 7.12E+00 1.685E-01 2.19E+00 1.69E+00 2.15E+01 1.60E-02 5.27E-07 5.12E+00 5.12E+00 5.12E+00 5.12E+00 5.12E+00 5.12E+00 5.12E+01 2.14E+01 2.14E+01 2.14E+01	based TRV (mg/kg-d) Conc. in soil (mg/kg) 3.79E+00 2.24E+00 7.04E-01 4.84E+00 1.48E+01 2.62E+02 5.29E-01 9.76E-01 6.85E-01 2.00E+02 7.12E+00 5.60E+01 1.16E+01 1.06E+02 6.85E-01 4.43E-01 2.19E+00 2.40E+02 1.69E+00 3.89E+01 2.15E+01 1.78E+02 1.60E-02 1.38E-01 5.27E-07 2.76E-05 5.12E+00 2.99E-01 5.12E+00 2.08E-01 5.12E+00 6.34E-01 5.12E+00 6.34E-01 2.14E+01 2.04E+00 2.14E+01 5.17E+00 5.27E-01 1.41E+00 2.14E+01 2.97E+01	based TRV (mg/kg-d) Conc. in soil (mg/kg) BTF 3.79E+00 2.24E+00 (a) 7.04E-01 4.84E+00 (a) 1.48E+01 2.62E+02 (a) 5.29E-01 9.76E-01 (a) 6.85E-01 2.00E+02 (a) 7.12E+00 5.60E+01 (a) 1.16E+01 1.06E+02 (a) 6.85E-01 4.43E-01 7.5E-06 2.19E+00 2.40E+02 (a) 6.85E-01 4.43E-01 7.5E-06 2.19E+00 2.40E+02 (a) 1.69E+00 3.89E+01 1.2E-02 2.15E+01 1.78E+02 (a) 1.60E-02 1.38E-01 5.8E-05 5.27E-07 2.76E-05 7.8E-05 5.12E+00 2.99E-01 0.0E+00 5.12E+00 3.13E-01 0.0E+00 5.12E+00 3.13E-01 0.0E+00 5.12E+00 6.34E-01 0.0E+00 5.12E+00 6.34E-01 0.0E+00 2.14E+01 2	NOAEL-based TRV (mg/kg-d) Conc. in soil (mg/kg) BTF (mg/kg) mammals (mg/kg) 3.79E+00 2.24E+00 (a) 3.5E-02 7.04E-01 4.84E+00 (a) 6.9E-02 1.48E+01 2.62E+02 (a) 4.8E-01 5.29E-01 9.76E-01 (a) 2.8E-01 6.85E-01 2.00E+02 (a) 1.7E+01 7.12E+00 5.60E+01 (a) 2.4E+00 1.16E+01 1.06E+02 (a) 8.1E+00 6.85E-01 4.43E-01 7.5E-06 3.3E-06 2.19E+00 2.40E+02 (a) 1.0E+01 1.69E+00 3.89E+01 1.2E-02 4.8E-01 2.15E+01 1.78E+02 (a) 1.1E+02 1.60E-02 1.38E-01 5.8E-05 8.0E-06 5.27E-07 2.76E-05 7.8E-05 2.2E-09 5.12E+00 2.08E-01 0.0E+00 0.0E+00 5.12E+00 2.08E-01 0.0E+00 0.0E+00 5.12E+00 6.34E-01 0.0E+00 0.0E	NOAEL-based TRV (mg/kg) (mg/kg	NOAEL- based TRV Conc. in soil (mg/kg) BTF (mg/kg) (mg/kg-d) (mg/kg-d	NOAEL- based TRV Conc. in soil (mg/kg) BTF (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg) (mg/kg-d) (mg/	NOAEL- based TRV Conc. in soil (mg/kg-d) (mg/kg) BTF (mg/kg) (mg/kg-d) (mg/kg				

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Attachment F-1 Table 1B - Evaluation of Potential Risk to Red Fox Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Red Fox

Pathway: Soil ingestion

Consumption of small mammals

Parameter	Value	
Small Mammal Ingestion Rate (kg ww/d)	0.510	Assume 100% of diet.
Total Dietary Intake (kg ww/d)	0.510	USEPA (1993).
Soil Ingestion Rate (kg dw/day)	0.01122	Calc. from Beyer (1994)
Body Weight (kg)	4.54	USEPA (1993). Mean of male and female BWs in spring and fall.
Sm. Mammal Dry wt./wet wt. CF	0.3	USEPA (1999)
Home range (ha)	1038	USEPA (1993)
Area Use Factor	0.00178	1.85. ha exposure area

	ĺ	Surface Soil									
				Conc. in		Intake from					
	LOAEL-			small	Intake	small					
	based TRV	Conc. in soil		mammals	from soil	mammals		TQ-small			
Constituent	(mg/kg-d)	(mg/kg)	BTF	(mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	mammals	Total TQ		
Antimony	7.86E-01	2.24E+00	(a)	3.5E-02	9.86E-06	2.09E-06	1.25E-05	2.66E-06	1.5E-05		
Arsenic	1.30E+00	4.84E+00	(a)	6.9E-02	2.13E-05	4.13E-06	1.64E-05	3.18E-06	2.0E-05		
Barium	2.36E+01	2.62E+02	(a)	4.8E-01	1.16E-03	2.88E-05	4.91E-05	1.22E-06	5.0E-05		
Cadmium	1.97E+00	9.76E-01	(a)	2.8E-01	4.30E-06	1.69E-05	2.18E-06	8.58E-06	1.1E-05		
Chromium	1.66E+01	2.00E+02	(a)	1.7E+01	8.82E-04	1.01E-03	5.32E-05	6.06E-05	1.1E-04		
Copper	2.36E+01	5.60E+01	(a)	2.4E+00	2.47E-04	1.44E-04	1.05E-05	6.12E-06	1.7E-05		
Lead	5.31E+01	1.06E+02	(a)	8.1E+00	4.69E-04	4.89E-04	8.83E-06	9.21E-06	1.8E-05		
Mercury	6.85E+00	4.43E-01	7.5E-06	3.3E-06	1.95E-06	6.67E-10	2.85E-07	9.74E-11	2.9E-07		
Nickel	4.21E+00	2.40E+02	(a)	1.0E+01	1.06E-03	6.03E-04	2.51E-04	1.43E-04	3.9E-04		
Vanadium	2.69E+00	3.89E+01	1.2E-02	4.8E-01	1.71E-04	2.87E-05	6.36E-05	1.07E-05	7.4E-05		
Zinc	8.48E+01	1.78E+02	(a)	1.1E+02	7.83E-04	6.80E-03	9.22E-06	8.01E-05	8.9E-05		
Aroclor-1260	1.60E-01	1.38E-01	5.8E-05	8.0E-06	6.07E-07	1.61E-09	3.79E-06	1.00E-08	3.8E-06		
2,3,7,8-TCDD TEQ (mammalian)	5.27E-06	2.76E-05	7.8E-05	2.2E-09	1.22E-10	4.31E-13	2.31E-05	8.19E-08	2.3E-05		
Benzo(b)fluoranthene	1.09E+01	2.99E-01	0.0E+00	0.0E+00	1.32E-06	0.00E+00	1.20E-07	0.00E+00	1.2E-07		
Benzo(g,h,i)perylene	1.09E+01	2.08E-01	0.0E+00	0.0E+00	9.17E-07	0.00E+00	8.38E-08	0.00E+00	8.4E-08		
Benzo(k)fluoranthene	1.09E+01	3.13E-01	0.0E+00	0.0E+00	1.38E-06	0.00E+00	1.26E-07	0.00E+00	1.3E-07		
Dibenz(a,h)anthracene	1.09E+01	2.08E-01	0.0E+00	0.0E+00	9.17E-07	0.00E+00	8.38E-08	0.00E+00	8.4E-08		
Pyrene	1.09E+01	6.34E-01	0.0E+00	0.0E+00	2.79E-06	0.00E+00	2.55E-07	0.00E+00	2.5E-07		
C5-C8 Aliphatics	2.14E+02	2.04E+00	1.0E+00	2.0E+00	9.00E-06	1.23E-04	4.19E-08	5.72E-07	6.1E-07		
C9-C10 Aromatics	2.14E+02	5.17E+00	1.0E+00	5.2E+00	2.28E-05	3.11E-04	1.07E-07	1.45E-06	1.6E-06		
C9-C18 Aliphatics	5.27E+02	1.41E+00	1.0E+00	1.4E+00	6.21E-06	8.46E-05	1.18E-08	1.61E-07	1.7E-07		
C11-C22 Aromatics	2.14E+02	2.97E+01	1.0E+00	3.0E+01	1.31E-04	1.78E-03	6.11E-07	8.33E-06	8.9E-06		
C19-C36 Aliphatics	1.05E+03	5.14E+01	1.0E+00	5.1E+01	2.26E-04	3.09E-03	2.15E-07	2.93E-06	3.1E-06		

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Table 2A - Evaluation of Potential Risk to Short-tailed Shrew Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA

Short-tailed Shrew Receptor: Pathway: Soil ingestion

Consumption of invertebrates

Parameter Value Invert Ingestion Rate (kg ww/d)
Total Dietary Intake (kg ww/d) Assumed to be 100% of diet 0.0080 0.0080 USEPA (1993) A Male & Female Soil Ingestion Rate (kg dw/day) 0.00012 Calc. from Beyer (1994) Body Weight (kg) 0.0168 USEPA (1993) Invert Dry wt./wet wt. CF 0.373 Site-specific data. Home range (ha) 0.39 USEPA (1993) Area Use Factor 1.85. ha exposure area

			Surface Soil								
						Intake					
	NOAEL-				Intake	from					
	based TRV	Conc. in soil		Conc. in	from soil	invert					
Constituent	(mg/kg-d)	(mg/kg)	BTF	invert (mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	TQ-invert	Total TQ		
Antimony	1.54E+01	2.24E+00	(a)	2.6E-01	1.66E-02	4.59E-02	1.08E-03	2.98E-03	4.1E-03		
Arsenic	2.85E+00	4.84E+00	(a)	1.3E+00	3.60E-02	2.29E-01	1.26E-02	8.04E-02	9.3E-02		
Barium	5.99E+01	2.62E+02	9.1E-02	2.4E+01	1.95E+00	4.21E+00	3.26E-02	7.04E-02	1.0E-01		
Cadmium	2.15E+00	9.76E-01	(a)	8.1E+00	7.25E-03	1.43E+00	3.38E-03	6.66E-01	6.7E-01		
Chromium	2.78E+00	2.00E+02	(a)	1.6E+01	1.49E+00	2.86E+00	5.36E-01	1.03E+00	1.6E+00		
Copper	2.88E+01	5.60E+01	(a)	1.5E+01	4.16E-01	2.61E+00	1.44E-02	9.05E-02	1.0E-01		
Lead	4.71E+01	1.06E+02	(a)	3.6E+01	7.91E-01	6.39E+00	1.68E-02	1.36E-01	1.5E-01		
Mercury	2.78E+00	4.43E-01	(a)	4.5E-01	3.30E-03	7.94E-02	1.19E-03	2.86E-02	3.0E-02		
Nickel	8.90E+00	2.40E+02	(a)	9.7E+00	1.78E+00	1.71E+00	2.00E-01	1.92E-01	3.9E-01		
Vanadium	6.85E+00	3.89E+01	(a)	5.2E+00	2.89E-01	9.17E-01	4.22E-02	1.34E-01	1.8E-01		
Zinc	8.71E+01	1.78E+02	(a)	1.7E+02	1.32E+00	2.96E+01	1.52E-02	3.40E-01	3.6E-01		
Aroclor-1260	6.50E-02	1.38E-01	1.1E+00	1.6E-01	1.02E-03	7.36E-02	1.58E-02	1.13E+00	1.1E+00		
2,3,7,8-TCDD TEQ (mammalian)	2.14E-06	2.76E-05	1.5E+00	4.0E-05	2.05E-07	1.89E-05	9.60E-02	8.86E+00	9.0E+00		
Benzo(b)fluoranthene	2.08E+01	2.99E-01	2.6E+00	7.8E-01	2.22E-03	1.37E-01	1.07E-04	6.61E-03	6.7E-03		
Benzo(g,h,i)perylene	2.08E+01	2.08E-01	2.9E+00	6.1E-01	1.55E-03	1.08E-01	7.45E-05	5.20E-03	5.3E-03		
Benzo(k)fluoranthene	2.08E+01	3.13E-01	2.6E+00	8.1E-01	2.33E-03	1.44E-01	1.12E-04	6.91E-03	7.0E-03		
Dibenz(a,h)anthracene	2.08E+01	2.08E-01	2.3E+00	4.8E-01	1.55E-03	8.48E-02	7.45E-05	4.08E-03	4.2E-03		
Pyrene	2.08E+01	6.34E-01	1.8E+00	1.1E+00	4.71E-03	1.96E-01	2.27E-04	9.42E-03	9.6E-03		
C5-C8 Aliphatics	8.70E+01	2.04E+00	1.0E+00	2.0E+00	1.52E-02	3.60E-01	1.75E-04	4.14E-03	4.3E-03		
C9-C10 Aromatics	8.67E+01	5.17E+00	1.0E+00	5.2E+00	3.85E-02	9.13E-01	4.44E-04	1.05E-02	1.1E-02		
C9-C18 Aliphatics	2.14E+02	1.41E+00	1.0E+00	1.4E+00	1.05E-02	2.48E-01	4.90E-05	1.16E-03	1.2E-03		
C11-C22 Aromatics	8.67E+01	2.97E+01	1.0E+00	3.0E+01	2.21E-01	5.23E+00	2.54E-03	6.04E-02	6.3E-02		
C19-C36 Aliphatics	4.27E+02	5.14E+01	1.0E+00	5.1E+01	3.82E-01	9.06E+00	8.94E-04	2.12E-02	2.2E-02		

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) Earthworm data collected at the site was used in place of BTF. Earthworm data used were from Sample DD-BO-001

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382.

Area Use Factor

Table 2B - Evaluation of Potential Risk to Short-tailed Shrew Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Short-tailed Shrew Pathway: Soil ingestion

Consumption of invertebrates

 Parameter
 Value

 Invert Ingestion Rate (kg ww/d)
 0.0080

 Total Dietary Intake (kg ww/d)
 0.0080

 Soil Ingestion Rate (kg dw/day)
 0.00012

 Body Weight (kg)
 0.0168

 Invert Dry wt./wet wt. CF
 0.373

 Home range (ha)
 0.39

Assumed to be 100% of diet USEPA (1993) A Male & Female

Calc. from Beyer (1994) USEPA (1993) Site-specific data. USEPA (1993) 1.85. ha exposure area

					Surface	Soil			
	LOAEL-				Intake	Intake from			
	based TRV	Conc. in soil		Conc. in	from soil	invert			
Constituent	(mg/kg-d)	(mg/kg)	BTF	invert (mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	TQ-invert	Total TQ
Antimony	3.19E+00	2.24E+00	(a)	2.6E-01	1.66E-02	4.59E-02	5.22E-03	1.44E-02	2.0E-02
Arsenic	5.26E+00	4.84E+00	(a)	1.3E+00	3.60E-02	2.29E-01	6.84E-03	4.36E-02	5.0E-02
Barium	9.55E+01	2.62E+02	9.1E-02	2.4E+01	1.95E+00	4.21E+00	2.04E-02	4.41E-02	6.5E-02
Cadmium	7.98E+00	9.76E-01	(a)	8.1E+00	7.25E-03	1.43E+00	9.09E-04	1.79E-01	1.8E-01
Chromium	6.72E+01	2.00E+02	(a)	1.6E+01	1.49E+00	2.86E+00	2.21E-02	4.25E-02	6.5E-02
Copper	9.56E+01	5.60E+01	(a)	1.5E+01	4.16E-01	2.61E+00	4.35E-03	2.73E-02	3.2E-02
Lead	2.15E+02	1.06E+02	(a)	3.6E+01	7.91E-01	6.39E+00	3.67E-03	2.96E-02	3.3E-02
Mercury	2.78E+01	4.43E-01	(a)	4.5E-01	3.30E-03	7.94E-02	1.19E-04	2.86E-03	3.0E-03
Nickel	1.71E+01	2.40E+02	(a)	9.7E+00	1.78E+00	1.71E+00	1.04E-01	1.00E-01	2.0E-01
Vanadium	1.09E+01	3.89E+01	(a)	5.2E+00	2.89E-01	9.17E-01	2.65E-02	8.41E-02	1.1E-01
Zinc	3.44E+02	1.78E+02	(a)	1.7E+02	1.32E+00	2.96E+01	3.84E-03	8.61E-02	9.0E-02
Aroclor-1260	6.50E-01	1.38E-01	1.1E+00	1.6E-01	1.02E-03	7.36E-02	1.58E-03	1.13E-01	1.1E-01
2,3,7,8-TCDD TEQ (mammalian)	2.14E-05	2.76E-05	1.5E+00	4.0E-05	2.05E-07	1.89E-05	9.60E-03	8.86E-01	9.0E-01
Benzo(b)fluoranthene	4.44E+01	2.99E-01	2.6E+00	7.8E-01	2.22E-03	1.37E-01	5.01E-05	3.09E-03	3.1E-03
Benzo(g,h,i)perylene	4.44E+01	2.08E-01	2.9E+00	6.1E-01	1.55E-03	1.08E-01	3.49E-05	2.43E-03	2.5E-03
Benzo(k)fluoranthene	4.44E+01	3.13E-01	2.6E+00	8.1E-01	2.33E-03	1.44E-01	5.24E-05	3.24E-03	3.3E-03
Dibenz(a,h)anthracene	4.44E+01	2.08E-01	2.3E+00	4.8E-01	1.55E-03	8.48E-02	3.49E-05	1.91E-03	1.9E-03
Pyrene	4.44E+01	6.34E-01	1.8E+00	1.1E+00	4.71E-03	1.96E-01	1.06E-04	4.41E-03	4.5E-03
C5-C8 Aliphatics	8.70E+02	2.04E+00	1.0E+00	2.0E+00	1.52E-02	3.60E-01	1.75E-05	4.14E-04	4.3E-04
C9-C10 Aromatics	8.67E+02	5.17E+00	1.0E+00	5.2E+00	3.85E-02	9.13E-01	4.44E-05	1.05E-03	1.1E-03
C9-C18 Aliphatics	2.14E+03	1.41E+00	1.0E+00	1.4E+00	1.05E-02	2.48E-01	4.90E-06	1.16E-04	1.2E-04
C11-C22 Aromatics	8.67E+02	2.97E+01	1.0E+00	3.0E+01	2.21E-01	5.23E+00	2.54E-04	6.04E-03	6.3E-03
C19-C36 Aliphatics	4.27E+03	5.14E+01	1.0E+00	5.1E+01	3.82E-01	9.06E+00	8.94E-05	2.12E-03	2.2E-03

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) Earthworm data collected at the site was used in place of BTF. Earthworm data used was from Sample DD-BO-001

References:

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382.

Attachment F-1 Table 3A - Evaluation of Potential Risk to Meadow Vole Bird Machine Company - Walpole, MA **Human Health and Environmental Risk Charaterization**

Site: DDA Meadow Vole Receptor: Pathway: Soil ingestion Consumption of vegetation

Parameter	Value	
Vegetation Ingestion Rate (kg ww/d)	0.0112	Assume 100% of diet.
Total Dietary Intake (kg ww/d)	0.0112	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.00021	Calc. from Beyer (1994)
Body Weight (kg)	0.0373	USEPA (1993) avg of A Male & Female all year
Veg Dry wt./wet wt. CF	0.120	USEPA (1999)
Home range (ha)	0.06	USEPA (1993) avg of A Male & Female grassy meadow MA
Area Use Factor	1	1.85. ha exposure area

					Surface	e Soil			
Constituent	NOAEL- based TRV (mg/kg-d)	Conc. in soil (mg/kg)	BTF	Conc. in vegetation (mg/kg)	Intake from soil (mg/kg-d)	Intake from vegetation (mg/kg-d)	TQ-soil	TQ- vegetation	Total TQ
Antimony	1.26E+01	2.24E+00	(a)	8.4E-02	1.27E-02	3.02E-03	1.01E-03	2.40E-04	1.2E-03
Arsenic	2.34E+00	4.84E+00	3.8E-02	1.8E-01	2.74E-02	6.54E-03	1.17E-02	2.80E-03	1.5E-02
Barium	4.90E+01	2.62E+02	1.6E-01	4.1E+01	1.49E+00	1.47E+00	3.03E-02	3.01E-02	6.0E-02
Cadmium	1.76E+00	9.76E-01	(a)	6.1E-01	5.52E-03	2.21E-02	3.14E-03	1.26E-02	1.6E-02
Chromium	2.28E+00	2.00E+02	4.1E-02	8.2E+00	1.13E+00	2.96E-01	4.98E-01	1.30E-01	6.3E-01
Copper	2.36E+01	5.60E+01	(a)	9.7E+00	3.17E-01	3.50E-01	1.34E-02	1.48E-02	2.8E-02
Lead	3.86E+01	1.06E+02	(a)	3.6E+00	6.03E-01	1.31E-01	1.56E-02	3.39E-03	1.9E-02
Mercury	2.28E+00	4.43E-01	4.3E-02	1.9E-02	2.51E-03	5.72E-03	1.10E-03	2.51E-03	3.6E-03
Nickel	7.29E+00	2.40E+02	(a)	6.5E+00	1.36E+00	2.35E-01	1.86E-01	3.22E-02	2.2E-01
Vanadium	5.61E+00	3.89E+01	4.9E-03	1.9E-01	2.20E-01	6.78E-03	3.92E-02	1.21E-03	4.0E-02
Zinc	7.14E+01	1.78E+02	(a)	8.5E+01	1.01E+00	3.07E+00	1.41E-02	4.29E-02	5.7E-02
Aroclor-1260	5.32E-02	1.38E-01	1.0E-02	1.4E-03	7.80E-04	4.13E-04	1.47E-02	7.77E-03	2.2E-02
2,3,7,8-TCDD TEQ (mammalian)	1.75E-06	2.76E-05	5.6E-03	1.5E-07	1.56E-07	5.56E-09	8.92E-02	3.18E-03	9.2E-02
Benzo(b)fluoranthene	1.70E+01	2.99E-01	3.1E-01	9.3E-02	1.69E-03	3.34E-03	9.95E-05	1.96E-04	3.0E-04
Benzo(g,h,i)perylene	1.70E+01	2.08E-01	(a)	6.2E-02	1.18E-03	2.22E-03	6.92E-05	1.30E-04	2.0E-04
Benzo(k)fluoranthene	1.70E+01	3.13E-01	(a)	4.3E-02	1.77E-03	1.53E-03	1.04E-04	9.01E-05	1.9E-04
Dibenz(a,h)anthracene	1.70E+01	2.08E-01	1.3E-01	2.7E-02	1.18E-03	9.74E-04	6.92E-05	5.73E-05	1.3E-04
Pyrene	1.70E+01	6.34E-01	7.2E-01	4.6E-01	3.59E-03	1.64E-02	2.11E-04	9.65E-04	1.2E-03
C5-C8 Aliphatics	7.12E+01	2.04E+00	1.0E+00	2.0E+00	1.16E-02	7.35E-02	1.62E-04	1.03E-03	1.2E-03
C9-C10 Aromatics	7.10E+01	5.17E+00	1.0E+00	5.2E+00	2.93E-02	1.86E-01	4.12E-04	2.62E-03	3.0E-03
C9-C18 Aliphatics	1.75E+02	1.41E+00	1.0E+00	1.4E+00	7.97E-03	5.07E-02	4.55E-05	2.90E-04	3.4E-04
C11-C22 Aromatics	7.10E+01	2.97E+01	1.0E+00	3.0E+01	1.68E-01	1.07E+00	2.36E-03	1.50E-02	1.7E-02
C19-C36 Aliphatics	3.50E+02	5.14E+01	1.0E+00	5.1E+01	2.91E-01	1.85E+00	8.30E-04	5.28E-03	6.1E-03

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References:

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Attachment F-1
Table 3B - Evaluation of Potential Risk to Meadow Vole
Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Meadow Vole

Pathway: Soil ingestion

Consumption of vegetation

Parameter	Value	
Vegetation Ingestion Rate (kg ww/d)	0.0112	Assume 100% of diet.
Total Dietary Intake (kg ww/d)	0.0112	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.00021	Calc. from Beyer (1994)
Body Weight (kg)	0.0373	USEPA (1993) avg of A Male & Female all year
Veg Dry wt./wet wt. CF	0.120	USEPA (1999)
Home range (ha)	0.06	USEPA (1993) avg of A Male & Female grassy meadow MA
Area Use Factor	1	1.85. ha exposure area

					Surface	Soil			
Constituent	LOAEL- based TRV (mg/kg-d)	Conc. in soil (mg/kg)	BTF	Conc. in vegetation (mg/kg)	Intake from soil (mg/kg-d)	Intake from vegetation (mg/kg-d)	TQ-soil	TQ- vegetation	Total TQ
Antimony	2.61E+00	2.24E+00	(a)	8.4E-02	1.27E-02	3.02E-03	4.85E-03	1.16E-03	6.0E-03
Arsenic	4.31E+00	4.84E+00	3.8E-02	1.8E-01	2.74E-02	6.54E-03	6.36E-03	1.52E-03	7.9E-03
Barium	7.83E+01	2.62E+02	1.6E-01	4.1E+01	1.49E+00	1.47E+00	1.90E-02	1.88E-02	3.8E-02
Cadmium	6.54E+00	9.76E-01	(a)	6.1E-01	5.52E-03	2.21E-02	8.45E-04	3.38E-03	4.2E-03
Chromium	5.51E+01	2.00E+02	4.1E-02	8.2E+00	1.13E+00	2.96E-01	2.06E-02	5.37E-03	2.6E-02
Copper	7.83E+01	5.60E+01	(a)	9.7E+00	3.17E-01	3.50E-01	4.05E-03	4.47E-03	8.5E-03
Lead	1.77E+02	1.06E+02	(a)	3.6E+00	6.03E-01	1.31E-01	3.41E-03	7.41E-04	4.2E-03
Mercury	2.28E+01	4.43E-01	4.3E-02	1.9E-02	2.51E-03	5.72E-03	1.10E-04	2.51E-04	3.6E-04
Nickel	1.40E+01	2.40E+02	(a)	6.5E+00	1.36E+00	2.35E-01	9.70E-02	1.68E-02	1.1E-01
Vanadium	8.94E+00	3.89E+01	4.9E-03	1.9E-01	2.20E-01	6.78E-03	2.46E-02	7.59E-04	2.5E-02
Zinc	2.82E+02	1.78E+02	(a)	8.5E+01	1.01E+00	3.07E+00	3.57E-03	1.09E-02	1.4E-02
Aroclor-1260	5.32E-01	1.38E-01	1.0E-02	1.4E-03	7.80E-04	4.13E-04	1.47E-03	7.77E-04	2.2E-03
2,3,7,8-TCDD TEQ (mammalian)	1.75E-05	2.76E-05	5.6E-03	1.5E-07	1.56E-07	5.56E-09	8.92E-03	3.18E-04	9.2E-03
Benzo(b)fluoranthene	3.64E+01	2.99E-01	3.1E-01	9.3E-02	1.69E-03	3.34E-03	4.66E-05	9.18E-05	1.4E-04
Benzo(g,h,i)perylene	3.64E+01	2.08E-01	(a)	6.2E-02	1.18E-03	2.22E-03	3.24E-05	6.09E-05	9.3E-05
Benzo(k)fluoranthene	3.64E+01	3.13E-01	(a)	4.3E-02	1.77E-03	1.53E-03	4.87E-05	4.22E-05	9.1E-05
Dibenz(a,h)anthracene	3.64E+01	2.08E-01	1.3E-01	2.7E-02	1.18E-03	9.74E-04	3.24E-05	2.68E-05	5.9E-05
Pyrene	3.64E+01	6.34E-01	7.2E-01	4.6E-01	3.59E-03	1.64E-02	9.86E-05	4.52E-04	5.5E-04
C5-C8 Aliphatics	7.12E+02	2.04E+00	1.0E+00	2.0E+00	1.16E-02	7.35E-02	1.62E-05	1.03E-04	1.2E-04
C9-C10 Aromatics	7.10E+02	5.17E+00	1.0E+00	5.2E+00	2.93E-02	1.86E-01	4.12E-05	2.62E-04	3.0E-04
C9-C18 Aliphatics	1.75E+03	1.41E+00	1.0E+00	1.4E+00	7.97E-03	5.07E-02	4.55E-06	2.90E-05	3.4E-05
C11-C22 Aromatics	7.10E+02	2.97E+01	1.0E+00	3.0E+01	1.68E-01	1.07E+00	2.36E-04	1.50E-03	1.7E-03
C19-C36 Aliphatics	3.50E+03	5.14E+01	1.0E+00	5.1E+01	2.91E-01	1.85E+00	8.30E-05	5.28E-04	6.1E-04

Notes:

kg ww/d - kilograms wet weight per day

kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Table 4A - Evaluation of Potential Risk to Red-tailed Hawk Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA Red-tailed Hawk Receptor: Pathway: Soil ingestion

Consumption of small mammals

Parameter	Value	
Small Mammal Ingestion Rate (kg ww/d)	0.119	Assumed 100% of diet
Total Dietary Intake (kg ww/d)	0.119	USEPA (1993) average of A Male & Female winter
Soil Ingestion Rate (kg dw/day)	0.00973	Calc. from Beyer (1994)
Body Weight (kg)	1.13	USEPA (1993), average of six adults
Sm. Mammal Dry wt./wet wt. CF	0.3	USEPA (1999)
Home range (ha)	60	USEPA (1993) A Male & Female spring
Area Use Factor	0.03083	1.85. ha exposure area

					Surface	Soil			
				Conc. in		Intake from			
	NOAEL-			small	Intake	small			
	based TRV	Conc. in soil		mammals	from soil	mammals		TQ-small	
Constituent	(mg/kg-d)	(mg/kg)	BTF	(mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	mammals	Total TQ
Antimony	1.00E+00	2.24E+00	(a)	3.5E-02	5.92E-04	3.39E-05	5.92E-04	3.39E-05	6.3E-04
Arsenic	3.70E+00	4.84E+00	(a)	6.9E-02	1.28E-03	6.68E-05	3.46E-04	1.80E-05	3.6E-04
Barium	2.08E+01	2.62E+02	(a)	4.8E-01	6.94E-02	4.66E-04	3.34E-03	2.24E-05	3.4E-03
Cadmium	1.47E+00	9.76E-01	(a)	2.8E-01	2.58E-04	2.73E-04	1.76E-04	1.86E-04	3.6E-04
Chromium	2.66E+00	2.00E+02	(a)	1.7E+01	5.30E-02	1.63E-02	1.99E-02	6.12E-03	2.6E-02
Copper	1.85E+01	5.60E+01	(a)	2.4E+00	1.48E-02	2.33E-03	8.01E-04	1.26E-04	9.3E-04
Lead	1.09E+01	1.06E+02	(a)	8.1E+00	2.82E-02	7.91E-03	2.58E-03	7.23E-04	3.3E-03
Mercury	4.50E-01	4.43E-01	7.5E-06	3.3E-06	1.17E-04	1.08E-08	2.61E-04	2.40E-08	2.6E-04
Nickel	6.71E+00	2.40E+02	(a)	1.0E+01	6.35E-02	9.75E-03	9.46E-03	1.45E-03	1.1E-02
Vanadium	1.19E+00	3.89E+01	1.2E-02	4.8E-01	1.03E-02	4.64E-04	8.67E-03	3.91E-04	9.1E-03
Zinc	6.61E+01	1.78E+02	(a)	1.1E+02	4.70E-02	1.10E-01	7.11E-04	1.66E-03	2.4E-03
Aroclor-1260	1.80E-01	1.38E-01	5.8E-05	8.0E-06	3.65E-05	2.60E-08	2.03E-04	1.45E-07	2.0E-04
2,3,7,8-TCDD TEQ (avian)	1.40E-05	4.92E-05	7.8E-05	3.8E-09	1.30E-08	1.24E-11	9.30E-04	8.88E-07	9.3E-04
Benzo(b)fluoranthene	1.01E+00	2.99E-01	0.0E+00	0.0E+00	7.92E-05	0.00E+00	7.84E-05	0.00E+00	7.8E-05
Benzo(g,h,i)perylene	1.01E+00	2.08E-01	0.0E+00	0.0E+00	5.51E-05	0.00E+00	5.45E-05	0.00E+00	5.5E-05
Benzo(k)fluoranthene	1.01E+00	3.13E-01	0.0E+00	0.0E+00	8.29E-05	0.00E+00	8.20E-05	0.00E+00	8.2E-05
Dibenz(a,h)anthracene	1.01E+00	2.08E-01	0.0E+00	0.0E+00	5.51E-05	0.00E+00	5.45E-05	0.00E+00	5.5E-05
Pyrene	1.11E+00	6.34E-01	0.0E+00	0.0E+00	1.68E-04	0.00E+00	1.51E-04	0.00E+00	1.5E-04
C5-C8 Aliphatics	4.00E+02	2.04E+00	1.0E+00	2.0E+00	5.40E-04	1.98E-03	1.35E-06	4.96E-06	6.3E-06
C9-C10 Aromatics	1.00E+04	5.17E+00	1.0E+00	5.2E+00	1.37E-03	5.03E-03	1.37E-07	5.03E-07	6.4E-07
C9-C18 Aliphatics	4.00E+02	1.41E+00	1.0E+00	1.4E+00	3.73E-04	1.37E-03	9.32E-07	3.42E-06	4.4E-06
C11-C22 Aromatics	1.00E+04	2.97E+01	1.0E+00	3.0E+01	7.85E-03	2.88E-02	7.85E-07	2.88E-06	3.7E-06
C19-C36 Aliphatics	4.00E+02	5.14E+01	1.0E+00	5.1E+01	1.36E-02	4.99E-02	3.40E-05	1.25E-04	1.6E-04

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Table 4B - Evaluation of Potential Risk to Red-tailed Hawk Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Red-tailed Hawk

Pathway: Soil ingestion

Consumption of small mammals

Parameter	Value	
Small Mammal Ingestion Rate (kg ww/d)	0.119	Assumed 100% of diet
Total Dietary Intake (kg ww/d)	0.119	USEPA (1993) average of A Male & Female winter
Soil Ingestion Rate (kg dw/day)	0.00973	Calc. from Beyer (1994)
Body Weight (kg)	1.13	USEPA (1993), average of six adults
Sm. Mammal Dry wt./wet wt. CF	0.3	USEPA (1999)
Home range (ha)	60	USEPA (1993) A Male & Female spring
Area Use Factor	0.03083	1.85. ha exposure area

		Surface Soil									
				Conc. in		Intake from					
	LOAEL-			small	Intake	small					
	based TRV	Conc. in soil		mammals	from soil	mammals		TQ-small			
Constituent	(mg/kg-d)	(mg/kg)	BTF	(mg/kg)	(mg/kg-d)	(mg/kg-d)	TQ-soil	mammals	Total TQ		
Antimony	5.00E+00	2.24E+00	(a)	3.5E-02	5.92E-04	3.39E-05	1.18E-04	6.77E-06	1.3E-04		
Arsenic	4.51E+00	4.84E+00	(a)	6.9E-02	1.28E-03	6.68E-05	2.84E-04	1.48E-05	3.0E-04		
Barium	4.17E+01	2.62E+02	(a)	4.8E-01	6.94E-02	4.66E-04	1.67E-03	1.12E-05	1.7E-03		
Cadmium	6.35E+00	9.76E-01	(a)	2.8E-01	2.58E-04	2.73E-04	4.07E-05	4.30E-05	8.4E-05		
Chromium	1.56E+01	2.00E+02	(a)	1.7E+01	5.30E-02	1.63E-02	3.39E-03	1.04E-03	4.4E-03		
Copper	3.49E+01	5.60E+01	(a)	2.4E+00	1.48E-02	2.33E-03	4.25E-04	6.69E-05	4.9E-04		
Lead	4.46E+01	1.06E+02	(a)	8.1E+00	2.82E-02	7.91E-03	6.31E-04	1.77E-04	8.1E-04		
Mercury	9.00E-01	4.43E-01	7.5E-06	3.3E-06	1.17E-04	1.08E-08	1.30E-04	1.20E-08	1.3E-04		
Nickel	1.86E+01	2.40E+02	(a)	1.0E+01	6.35E-02	9.75E-03	3.42E-03	5.25E-04	3.9E-03		
Vanadium	1.70E+00	3.89E+01	1.2E-02	4.8E-01	1.03E-02	4.64E-04	6.05E-03	2.73E-04	6.3E-03		
Zinc	1.71E+02	1.78E+02	(a)	1.1E+02	4.70E-02	1.10E-01	2.74E-04	6.41E-04	9.2E-04		
Aroclor-1260	1.80E+00	1.38E-01	5.8E-05	8.0E-06	3.65E-05	2.60E-08	2.03E-05	1.45E-08	2.0E-05		
2,3,7,8-TCDD TEQ (avian)	1.40E-04	4.92E-05	7.8E-05	3.8E-09	1.30E-08	1.24E-11	9.30E-05	8.88E-08	9.3E-05		
Benzo(b)fluoranthene	1.01E+01	2.99E-01	0.0E+00	0.0E+00	7.92E-05	0.00E+00	7.84E-06	0.00E+00	7.8E-06		
Benzo(g,h,i)perylene	1.01E+01	2.08E-01	0.0E+00	0.0E+00	5.51E-05	0.00E+00	5.45E-06	0.00E+00	5.5E-06		
Benzo(k)fluoranthene	1.01E+01	3.13E-01	0.0E+00	0.0E+00	8.29E-05	0.00E+00	8.20E-06	0.00E+00	8.2E-06		
Dibenz(a,h)anthracene	1.01E+01	2.08E-01	0.0E+00	0.0E+00	5.51E-05	0.00E+00	5.45E-06	0.00E+00	5.5E-06		
Pyrene	1.11E+01	6.34E-01	0.0E+00	0.0E+00	1.68E-04	0.00E+00	1.51E-05	0.00E+00	1.5E-05		
C5-C8 Aliphatics	4.00E+03	2.04E+00	1.0E+00	2.0E+00	5.40E-04	1.98E-03	1.35E-07	4.96E-07	6.3E-07		
C9-C10 Aromatics	1.00E+05	5.17E+00	1.0E+00	5.2E+00	1.37E-03	5.03E-03	1.37E-08	5.03E-08	6.4E-08		
C9-C18 Aliphatics	4.00E+03	1.41E+00	1.0E+00	1.4E+00	3.73E-04	1.37E-03	9.32E-08	3.42E-07	4.4E-07		
C11-C22 Aromatics	1.00E+05	2.97E+01	1.0E+00	3.0E+01	7.85E-03	2.88E-02	7.85E-08	2.88E-07	3.7E-07		
C19-C36 Aliphatics	4.00E+03	5.14E+01	1.0E+00	5.1E+01	1.36E-02	4.99E-02	3.40E-06	1.25E-05	1.6E-05		

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Table 5A - Evaluation of Potential Risk to American Woodcock Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

Site:

American Woodcock Receptor:

Pathway: Soil ingestion

Consumption of invertebrates

Parameter	Value	7
Invert Ingestion Rate (kg ww/d)	0.1371	Assumed 100% of diet
Total Dietary Intake (kg ww/d)	0.1371	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.0112	Calc. from Beyer (1994)
Body Weight (kg)	0.1780	USEPA (1993) avg of A Male & A Female in central MA
Invert Dry wt./wet wt. CF	0.373	Site-specific data.
Home range (ha)	3.80	USEPA (1993) avg of inactive A Male and brooding A Female
Area Use Factor	0.4868	1.85. ha exposure area

			Surface Soil						
	NOAEL-				Intake	Intake from			
	based TRV	Conc. in soil		Conc. in	from soil	invert			
Constituent	(mg/kg-d)	(mg/kg)	BTF	invert (mg/kg)		(mg/kg-d)	TQ-soil	TQ-invert	Total TQ
Antimony	1.00E+00	2.24E+00	(a)	2.6E-01	6.86E-02	3.64E-02	6.86E-02	3.64E-02	1.0E-01
Arsenic	3.70E+00	4.84E+00	(a)	1.3E+00	1.48E-01	1.82E-01	4.01E-02	4.91E-02	8.9E-02
Barium	2.08E+01	2.62E+02	9.1E-02	2.4E+01	8.04E+00	3.34E+00	3.87E-01	1.61E-01	5.5E-01
Cadmium	1.47E+00	9.76E-01	(a)	8.1E+00	2.99E-02	1.13E+00	2.04E-02	7.72E-01	7.9E-01
Chromium	2.66E+00	2.00E+02	(a)	1.6E+01	6.14E+00	2.27E+00	2.31E+00	8.52E-01	3.2E+00
Copper	1.85E+01	5.60E+01	(a)	1.5E+01	1.72E+00	2.07E+00	9.28E-02	1.12E-01	2.0E-01
Lead	1.09E+01	1.06E+02	(a)	3.6E+01	3.26E+00	5.06E+00	2.98E-01	4.63E-01	7.6E-01
Mercury	4.50E-01	4.43E-01	(a)	4.5E-01	1.36E-02	6.29E-02	3.02E-02	1.40E-01	1.7E-01
Nickel	6.71E+00	2.40E+02	(a)	9.7E+00	7.35E+00	1.36E+00	1.10E+00	2.02E-01	1.3E+00
Vanadium	1.19E+00	3.89E+01	(a)	5.2E+00	1.19E+00	7.27E-01	1.00E+00	6.13E-01	1.6E+00
Zinc	6.61E+01	1.78E+02	(a)	1.7E+02	5.44E+00	2.35E+01	8.24E-02	3.56E-01	4.4E-01
Aroclor-1260	1.80E-01	1.38E-01	1.1E+00	1.6E-01	4.22E-03	5.84E-02	2.35E-02	3.24E-01	3.5E-01
2,3,7,8-TCDD TEQ (avian)	1.40E-05	4.92E-05	1.5E+00	7.1E-05	1.51E-06	2.67E-05	1.08E-01	1.91E+00	2.0E+00
Benzo(b)fluoranthene	1.01E+00	2.99E-01	2.6E+00	7.8E-01	9.17E-03	1.09E-01	9.08E-03	1.08E-01	1.2E-01
Benzo(g,h,i)perylene	1.01E+00	2.08E-01	2.9E+00	6.1E-01	6.38E-03	8.56E-02	6.31E-03	8.47E-02	9.1E-02
Benzo(k)fluoranthene	1.01E+00	3.13E-01	2.6E+00	8.1E-01	9.60E-03	1.14E-01	9.50E-03	1.13E-01	1.2E-01
Dibenz(a,h)anthracene	1.01E+00	2.08E-01	2.3E+00	4.8E-01	6.38E-03	6.72E-02	6.31E-03	6.66E-02	7.3E-02
Pyrene	1.11E+00	6.34E-01	1.8E+00	1.1E+00	1.94E-02	1.55E-01	1.75E-02	1.40E-01	1.6E-01
C5-C8 Aliphatics	4.00E+02	2.04E+00	1.0E+00	2.0E+00	6.26E-02	2.85E-01	1.56E-04	7.14E-04	8.7E-04
C9-C10 Aromatics	1.00E+04	5.17E+00	1.0E+00	5.2E+00	1.59E-01	7.24E-01	1.59E-05	7.24E-05	8.8E-05
C9-C18 Aliphatics	4.00E+02	1.41E+00	1.0E+00	1.4E+00	4.32E-02	1.97E-01	1.08E-04	4.92E-04	6.0E-04
C11-C22 Aromatics	1.00E+04	2.97E+01	1.0E+00	3.0E+01	9.09E-01	4.15E+00	9.09E-05	4.15E-04	5.1E-04
C19-C36 Aliphatics	4.00E+02	5.14E+01	1.0E+00	5.1E+01	1.57E+00	7.18E+00	3.94E-03	1.80E-02	2.2E-02

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) Earthworm data collected at the site was used in place of BTF. Earthworm data used was from Sample DD-BO-001

References:

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382.

Table 5B - Evaluation of Potential Risk to American Woodcock Bird Machine Company - Walpole, MA

Human Health and Environmental Risk Characterization

DDA Site:

American Woodcock Receptor:

Pathway: Soil ingestion

Consumption of invertebrates

Parameter	Value	
Invert Ingestion Rate (kg ww/d)	0.1371	Assumed 100% of diet
Total Dietary Intake (kg ww/d)	0.1371	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.0112	Calc. from Beyer (1994)
Body Weight (kg)	0.1780	USEPA (1993) avg of A Male & A Female in central MA
Invert Dry wt./wet wt. CF	0.373	Site-specific data.
Home range (ha)	3.80	USEPA (1993) avg of inactive A Male and brooding A Female
Area Use Factor	0.4868	1.85. ha exposure area

		Surface Soil							
	LOAEL-				Intake	Intake from			
	based TRV	Conc. in soil		Conc. in	from soil	invert			
Constituent	(mg/kg-d)	(mg/kg)	BTF	invert (mg/kg)		(mg/kg-d)	TQ-soil	TQ-invert	Total TQ
Antimony	5.00E+00	2.24E+00	(a)	2.6E-01	6.86E-02	3.64E-02	1.37E-02	7.27E-03	2.1E-02
Arsenic	4.51E+00	4.84E+00	(a)	1.3E+00	1.48E-01	1.82E-01	3.29E-02	4.03E-02	7.3E-02
Barium	4.17E+01	2.62E+02	9.1E-02	2.4E+01	8.04E+00	3.34E+00	1.93E-01	8.01E-02	2.7E-01
Cadmium	6.35E+00	9.76E-01	(a)	8.1E+00	2.99E-02	1.13E+00	4.71E-03	1.78E-01	1.8E-01
Chromium	1.56E+01	2.00E+02	(a)	1.6E+01	6.14E+00	2.27E+00	3.93E-01	1.45E-01	5.4E-01
Copper	3.49E+01	5.60E+01	(a)	1.5E+01	1.72E+00	2.07E+00	4.92E-02	5.93E-02	1.1E-01
Lead	4.46E+01	1.06E+02	(a)	3.6E+01	3.26E+00	5.06E+00	7.31E-02	1.13E-01	1.9E-01
Mercury	9.00E-01	4.43E-01	(a)	4.5E-01	1.36E-02	6.29E-02	1.51E-02	6.99E-02	8.5E-02
Nickel	1.86E+01	2.40E+02	(a)	9.7E+00	7.35E+00	1.36E+00	3.96E-01	7.30E-02	4.7E-01
Vanadium	1.70E+00	3.89E+01	(a)	5.2E+00	1.19E+00	7.27E-01	7.00E-01	4.28E-01	1.1E+00
Zinc	1.71E+02	1.78E+02	(a)	1.7E+02	5.44E+00	2.35E+01	3.17E-02	1.37E-01	1.7E-01
Aroclor-1260	1.80E+00	1.38E-01	1.1E+00	1.6E-01	4.22E-03	5.84E-02	2.35E-03	3.24E-02	3.5E-02
2,3,7,8-TCDD TEQ (avian)	1.40E-04	4.92E-05	1.5E+00	7.1E-05	1.51E-06	2.67E-05	1.08E-02	1.91E-01	2.0E-01
Benzo(b)fluoranthene	1.01E+01	2.99E-01	2.6E+00	7.8E-01	9.17E-03	1.09E-01	9.08E-04	1.08E-02	1.2E-02
Benzo(g,h,i)perylene	1.01E+01	2.08E-01	2.9E+00	6.1E-01	6.38E-03	8.56E-02	6.31E-04	8.47E-03	9.1E-03
Benzo(k)fluoranthene	1.01E+01	3.13E-01	2.6E+00	8.1E-01	9.60E-03	1.14E-01	9.50E-04	1.13E-02	1.2E-02
Dibenz(a,h)anthracene	1.01E+01	2.08E-01	2.3E+00	4.8E-01	6.38E-03	6.72E-02	6.31E-04	6.66E-03	7.3E-03
Pyrene	1.11E+01	6.34E-01	1.8E+00	1.1E+00	1.94E-02	1.55E-01	1.75E-03	1.40E-02	1.6E-02
C5-C8 Aliphatics	4.00E+03	2.04E+00	1.0E+00	2.0E+00	6.26E-02	2.85E-01	1.56E-05	7.14E-05	8.7E-05
C9-C10 Aromatics	1.00E+05	5.17E+00	1.0E+00	5.2E+00	1.59E-01	7.24E-01	1.59E-06	7.24E-06	8.8E-06
C9-C18 Aliphatics	4.00E+03	1.41E+00	1.0E+00	1.4E+00	4.32E-02	1.97E-01	1.08E-05	4.92E-05	6.0E-05
C11-C22 Aromatics	1.00E+05	2.97E+01	1.0E+00	3.0E+01	9.09E-01	4.15E+00	9.09E-06	4.15E-05	5.1E-05
C19-C36 Aliphatics	4.00E+03	5.14E+01	1.0E+00	5.1E+01	1.57E+00	7.18E+00	3.94E-04	1.80E-03	2.2E-03

Notes:

kg ww/d - kilograms wet weight per day

kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) Earthworm data collected at the site was used in place of BTF. Earthworm data used was from Sample DD-BO-001

References:

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382.

Attachment F-1
Table 6A - Evaluation of Potential Risk to Quail
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Quail

Pathway: Soil ingestion

Consumption of vegetation

Parameter Value Vegetation Ingestion Rate (kg ww/d) Total Dietary Intake (kg ww/d) Assume 100% of diet. 0.0134 USEPA (1993) 0.0134 Soil Ingestion Rate (kg dw/day) 0.0002 Calc. from Beyer (1994) Body Weight (kg) 0.1736 USEPA (1993) avg of A Male and A Female Veg Dry wt./wet wt. CF USEPA (1999) 1.16E-02 0.120 Home range (ha) 9.98 USEPA (1993) avg of A Male and A Female Area Use Factor 0.185 1.85. ha exposure area

		Surface Soil							
Constituent	NOAEL- based TRV (mg/kg-d)	Conc. in soil (mg/kg)	BTF	Conc. in vegetation (mg/kg)	Intake from soil (mg/kg-d)	Intake from vegetation (mg/kg-d)	TQ-soil	TQ- vegetation	Total TQ
Antimony	1.00E+00	2.24E+00	(a)	8.4E-02	3.99E-04	1.44E-04	3.99E-04	1.44E-04	5.4E-04
Arsenic	3.70E+00	4.84E+00	3.8E-02	1.8E-01	8.63E-04	3.11E-04	2.33E-04	8.41E-05	3.2E-04
Barium	2.08E+01	2.62E+02	1.6E-01	4.1E+01	4.67E-02	7.01E-02	2.25E-03	3.37E-03	5.6E-03
Cadmium	1.47E+00	9.76E-01	(a)	6.1E-01	1.74E-04	1.05E-03	1.18E-04	7.16E-04	8.3E-04
Chromium	2.66E+00	2.00E+02	4.1E-02	8.2E+00	3.57E-02	1.41E-02	1.34E-02	5.29E-03	1.9E-02
Copper	1.85E+01	5.60E+01	(a)	9.7E+00	9.98E-03	1.66E-02	5.39E-04	9.00E-04	1.4E-03
Lead	1.09E+01	1.06E+02	(a)	3.6E+00	1.90E-02	6.23E-03	1.73E-03	5.69E-04	2.3E-03
Mercury	4.50E-01	4.43E-01	4.3E-02	1.9E-02	7.90E-05	2.72E-04	1.75E-04	6.05E-04	7.8E-04
Nickel	6.71E+00	2.40E+02	(a)	6.5E+00	4.27E-02	1.12E-02	6.37E-03	1.67E-03	8.0E-03
Vanadium	1.19E+00	3.89E+01	4.9E-03	1.9E-01	6.92E-03	3.23E-04	5.84E-03	2.72E-04	6.1E-03
Zinc	6.61E+01	1.78E+02	(a)	8.5E+01	3.16E-02	1.46E-01	4.79E-04	2.21E-03	2.7E-03
Aroclor-1260	1.80E-01	1.38E-01	1.0E-02	1.4E-03	2.45E-05	1.97E-05	1.36E-04	1.09E-04	2.5E-04
2,3,7,8-TCDD TEQ (avian)	1.40E-05	4.92E-05	5.6E-03	2.8E-07	8.76E-09	4.72E-10	6.26E-04	3.37E-05	6.6E-04
Benzo(b)fluoranthene	1.01E+00	2.99E-01	3.1E-01	9.3E-02	5.33E-05	1.59E-04	5.28E-05	1.57E-04	2.1E-04
Benzo(g,h,i)perylene	1.01E+00	2.08E-01	(a)	6.2E-02	3.71E-05	1.05E-04	3.67E-05	1.04E-04	1.4E-04
Benzo(k)fluoranthene	1.01E+00	3.13E-01	(a)	4.3E-02	5.58E-05	7.30E-05	5.52E-05	7.22E-05	1.3E-04
Dibenz(a,h)anthracene	1.01E+00	2.08E-01	1.3E-01	2.7E-02	3.71E-05	4.63E-05	3.67E-05	4.59E-05	8.3E-05
Pyrene	1.11E+00	6.34E-01	7.2E-01	4.6E-01	1.13E-04	7.81E-04	1.02E-04	7.04E-04	8.1E-04
C5-C8 Aliphatics	4.00E+02	2.04E+00	1.0E+00	2.0E+00	3.64E-04	3.50E-03	9.09E-07	8.74E-06	9.7E-06
C9-C10 Aromatics	1.00E+04	5.17E+00	1.0E+00	5.2E+00	9.22E-04	8.86E-03	9.22E-08	8.86E-07	9.8E-07
C9-C18 Aliphatics	4.00E+02	1.41E+00	1.0E+00	1.4E+00	2.51E-04	2.41E-03	6.27E-07	6.03E-06	6.7E-06
C11-C22 Aromatics	1.00E+04	2.97E+01	1.0E+00	3.0E+01	5.28E-03	5.08E-02	5.28E-07	5.08E-06	5.6E-06
C19-C36 Aliphatics	4.00E+02	5.14E+01	1.0E+00	5.1E+01	9.15E-03	8.80E-02	2.29E-05	2.20E-04	2.4E-04

Notes:

kg ww/d - kilograms wet weight per day kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor TQ - Task quotient

TRV - Toxicity reference value

NOAEL - No observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Attachment F-1
Table 6B - Evaluation of Potential Risk to Quail
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

Site: DDA

Receptor: Quail

Pathway: Soil ingestion

Consumption of vegetation

Parameter	Value	
Vegetation Ingestion Rate (kg ww/d)	0.0134	Assume 100% of diet.
Total Dietary Intake (kg ww/d)	0.0134	USEPA (1993)
Soil Ingestion Rate (kg dw/day)	0.0002	Calc. from Beyer (1994)
Body Weight (kg)	0.1736	USEPA (1993) avg of A Male and A Female
Veg Dry wt./wet wt. CF	0.120	USEPA (1999)
Home range (ha)	9.98	USEPA (1993) avg of A Male and A Female
Area Use Factor	0.185	1.85. ha exposure area

					Surface	e Soil			
Constituent	LOAEL- based TRV (mg/kg-d)	Conc. in soil (mg/kg)	BTF	Conc. in vegetation (mg/kg)	Intake from soil (mg/kg-d)	Intake from vegetation (mg/kg-d)	TQ-soil	TQ- vegetation	Total TQ
Antimony	5.00E+00	2.24E+00	(a)	8.4E-02	3.99E-04	1.44E-04	7.97E-05	2.88E-05	1.1E-04
Arsenic	4.51E+00	4.84E+00	3.8E-02	1.8E-01	8.63E-04	3.11E-04	1.91E-04	6.91E-05	2.6E-04
Barium	4.17E+01	2.62E+02	1.6E-01	4.1E+01	4.67E-02	7.01E-02	1.12E-03	1.68E-03	2.8E-03
Cadmium	6.35E+00	9.76E-01	(a)	6.1E-01	1.74E-04	1.05E-03	2.74E-05	1.66E-04	1.9E-04
Chromium	1.56E+01	2.00E+02	4.1E-02	8.2E+00	3.57E-02	1.41E-02	2.28E-03	9.00E-04	3.2E-03
Copper	3.49E+01	5.60E+01	(a)	9.7E+00	9.98E-03	1.66E-02	2.86E-04	4.77E-04	7.6E-04
Lead	4.46E+01	1.06E+02	(a)	3.6E+00	1.90E-02	6.23E-03	4.25E-04	1.40E-04	5.6E-04
Mercury	9.00E-01	4.43E-01	4.3E-02	1.9E-02	7.90E-05	2.72E-04	8.77E-05	3.02E-04	3.9E-04
Nickel	1.86E+01	2.40E+02	(a)	6.5E+00	4.27E-02	1.12E-02	2.30E-03	6.02E-04	2.9E-03
Vanadium	1.70E+00	3.89E+01	4.9E-03	1.9E-01	6.92E-03	3.23E-04	4.07E-03	1.90E-04	4.3E-03
Zinc	1.71E+02	1.78E+02	(a)	8.5E+01	3.16E-02	1.46E-01	1.85E-04	8.51E-04	1.0E-03
Aroclor-1260	1.80E+00	1.38E-01	1.0E-02	1.4E-03	2.45E-05	1.97E-05	1.36E-05	1.09E-05	2.5E-05
2,3,7,8-TCDD TEQ (avian)	1.40E-04	4.92E-05	5.6E-03	2.8E-07	8.76E-09	4.72E-10	6.26E-05	3.37E-06	6.6E-05
Benzo(b)fluoranthene	1.01E+01	2.99E-01	3.1E-01	9.3E-02	5.33E-05	1.59E-04	5.28E-06	1.57E-05	2.1E-05
Benzo(g,h,i)perylene	1.01E+01	2.08E-01	(a)	6.2E-02	3.71E-05	1.05E-04	3.67E-06	1.04E-05	1.4E-05
Benzo(k)fluoranthene	1.01E+01	3.13E-01	(a)	4.3E-02	5.58E-05	7.30E-05	5.52E-06	7.22E-06	1.3E-05
Dibenz(a,h)anthracene	1.01E+01	2.08E-01	1.3E-01	2.7E-02	3.71E-05	4.63E-05	3.67E-06	4.59E-06	8.3E-06
Pyrene	1.11E+01	6.34E-01	7.2E-01	4.6E-01	1.13E-04	7.81E-04	1.02E-05	7.04E-05	8.1E-05
C5-C8 Aliphatics	4.00E+03	2.04E+00	1.0E+00	2.0E+00	3.64E-04	3.50E-03	9.09E-08	8.74E-07	9.7E-07
C9-C10 Aromatics	1.00E+05	5.17E+00	1.0E+00	5.2E+00	9.22E-04	8.86E-03	9.22E-09	8.86E-08	9.8E-08
C9-C18 Aliphatics	4.00E+03	1.41E+00	1.0E+00	1.4E+00	2.51E-04	2.41E-03	6.27E-08	6.03E-07	6.7E-07
C11-C22 Aromatics	1.00E+05	2.97E+01	1.0E+00	3.0E+01	5.28E-03	5.08E-02	5.28E-08	5.08E-07	5.6E-07
C19-C36 Aliphatics	4.00E+03	5.14E+01	1.0E+00	5.1E+01	9.15E-03	8.80E-02	2.29E-06	2.20E-05	2.4E-05

Notes:

kg ww/d - kilograms wet weight per day

kg dw/d - kilograms dry weight per day

CF - conversion factor

mg/kg-day - milligrams per kilogram per day

BTF - Biotransfer factor

TQ - Task quotient

TRV - Toxicity reference value

LOAEL - Lowest observable adverse effect level

(a) BTF not available; ECO-SSL uptake equations are used

References

USEPA 1993. Wildlife Exposure Factors Handbook. EPA 600-R-93-187

Beyer W. Nelson. Connor Erin E. Gerould Sarah. 1994. Estimates of soil ingestion by wildlife. Journal of Wildlife Management. 58(2): 375-382. USEPA. 1999. Screening Level Ecological Risk Assessment Protocol for Hazardous Waste Combustion Facilities. EPA 530-D-99-001A.

Attachment F-2
Biotransfer Factors using Uptake Equations
Bird Machine Company - Walpole, MA
Human Health and Environmental Risk Characterization

	DDA Uptake Equations from EcoSSLs				
	Invertebrate	Vegetation	Small Mammal		
Constituent	BTFs	BTFs	BTFs		
Antimony		8.40E-02	3.49E-02		
Arsenic	7.35E-01		6.88E-02		
Barium			4.80E-01		
Benzo(a)anthracene					
Benzo(a)pyrene					
Benzo(g,h,i)perylene		6.15E-02			
Benzo(k)fluoranthene		4.26E-02			
Cadmium	8.12E+00	6.14E-01	2.81E-01		
Chromium			1.67E+01		
Copper		9.72E+00	2.40E+00		
Lead	3.48E+01	3.64E+00	8.15E+00		
Nickel		6.53E+00	1.00E+01		
Zinc	4.68E+02	8.52E+01	1.13E+02		

Notes:

All units are in dry weight soil.

Uptake equations are used in risk calculations whenever published BTFs (Table 4-4) are not available.



APPENDIX V REPRESENTATIVENESS AND DATA USABILITY WORKSHEET

A. Representativeness Evaluation (Specific to info Refer to Section 6.0 through 6.8.)	ormation/samples used to support the RAO.
A-1 Provide a succinct summary of the Conceptual Site Model (CSM) for the disposal site. Discussion should include: - Disposal site history - Geologic/hydrogeological setting - Contaminant Source(s) and Type(s) - Description of the volume/mass and types of contaminants released to the environment - Date/time period of release(s), if known - Release location, affected media, and horizontal and vertical extent of the contamination - Contaminant migration pathways - Mechanism/pathways and points of exposure by human and ecological receptors	
 A-2 Discuss use of Field/Screening Data in response action decision making, including: Contaminant of Concern screening/elimination Selection of sampling locations Comparison to laboratory results Comparison to visual/olfactory observations (Refer to Section 6.2) A-3 Discuss and justify sampling locations and depths collected in support of RAO regarding: For Class A or B RAOs Delineation of disposal site boundaries (horizontal and vertical) Elimination/control of OHM source(s) Characterization of Risk (Exposure Pathways/Receptors, Hot Spots, samples included in EPCs, Background) Achievement of No Significant Risk (NSR) For Class C RAOs Delineation of disposal site boundaries (horizontal 	() No Field/Screening Data were used to directly support this RAO.() Field/Screening Data were used, as follows:
and vertical) -Elimination/control of OHM source(s) -Characterization of Risk (Exposure Pathways/Receptors, Hot Spots, samples included in EPCs, Background) -Achievement of No Substantial Hazard (NSH) (Refer to Table1 and Section 6.3; A-3 and A-4 of the worksheet may be combined, as appropriate.)	

A. Representativeness Evaluation (Specific to info Refer to Section 6.0 through 6.8.)	ormation/samples used to support the RAO.
A-4 Discuss and justify the density, spatial distribution, collection methods, and handling (compositing, split sampling) of samples collected in support of RAO (in relation to the justification provided in A-3 for meeting the RAO requirements)	
(Refer to Table 1 and Section 6.4)	
A-5 Identify disposal site conditions, if any, that warrant the collection and analysis of temporal samples. For disposal sites that require monitoring over an extended time period, discuss and justify the number and time interval for sampling rounds conducted in support of the RAO for the following: For Class A or B RAOs - Delineation of disposal site boundaries (horizontal and vertical) - Characterization of Risk (Exposure Pathways/Receptors, Hot Spots, samples included in EPCs, Background) - Elimination/control of OHM source(s) - Achievement of No Significant Risk (NSR) For Class C RAOs - Delineation of disposal site boundaries (horizontal and vertical) - Characterization of Risk (Exposure Pathways/Receptors, Hot Spots, samples included in EPCs, Background) - Elimination/control of OHM source(s) - Achievement of No Substantial Hazard (NSH)	() Temporal sampling not warranted for this disposal site.
(Refer to Table 1 and Section 6.5)	
A-6 Field Completeness of Data: Discuss data gaps identified in sampling and analytical information used to support RAO and their significance.	
(Refer to Section 6.6)	
A-7 Identify any inconsistent information or uncertainty and justify disregarding such information or uncertainty (e.g., site assessment data inconsistent with historical information, field screening data/observations inconsistent with analytical data, use of data to support the RAO in spite of identified analytical or other deficiencies, etc.) in rendering the RAO Opinion.	
(Refer to Section 6.7)	

A. Representativeness Evaluation (Specific to info Refer to Section 6.0 through 6.8.)	ormation/samples used to support the RAO.
A-8 Where it is not otherwise apparent or discussed in previous sections, identify/discuss information generated during the course of response actions that was not used to support the RAO because it was determined to be unrepresentative or no longer representative of disposal site conditions.	
(Refer to Section 6.8)	

	B. <u>Data Usability Assessment</u> (Specific to samples used to support the RAO. Refer to Table 1, Section 7.0 through 7.3, and Appendices I, II, III and IV.)					
B-1 List all MCP activities that provided the analytical data reviewed in the course of conducting the Data Usability Assessment in support of the RAO. Include the media sampled and the month and year the data were acquired.	() Listed below. () Attached separately (provide attachment reference).					
B-2 Discuss appropriateness of selected analytical methods to quantitatively support the RAO.						
B-3 Discuss appropriateness of selected analytical methods' Reporting Limits (RL) to quantitatively support the RAO.	() All Reporting Limits were at or below applicable standards.					
B-4 Discuss laboratory performance criteria and data quality indicators used to assess overall Analytical Accuracy (continuing calibration, laboratory control spikes, etc.) and Analytical Precision (laboratory duplicates, laboratory control spike duplicates, etc.). For CAM data, see MCP Analytical Method Report Certification Form and Laboratory Case Narrative.	() Met all CAM requirements and performance standards without qualification.() If not, discuss data usability implications.					
B-5 Discuss performance criteria and data quality indicators used to assess overall Field Data Usability (sample preservation compliance, sample sub sampling/compositing, etc.).						
B-6 Discuss any data rejected pursuant to Appendix IV, Rejection Criteria – Analytical Data Usability Assessments.	() No data rejected pursuant to Appendix IV .					

C. Representativeness Evaluation and Data L Conclusions (Refer to Section 8.0)	Isability Assessment Summary and
Provide a summary declaration that the data set relied upon to support the RAO is:	
Scientifically valid and defensible, and of sufficient accuracy, precision and completeness; and	
Representative with regards to the spatial and temporal distribution of sampling points.	



Data Usability Assessment

Data in support of this RAO were reviewed to ensure applicable MCP guidelines and policies were followed in regards to data quality and data usability. The data usability assessment has both an analytical and a field component, each of which are provided in the following subsections. All analytical data were collected after August 2003 when the MADEP established "presumptive certainty" requirements as defined in "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data" (BWSC-CAM-VIIA, rev. 3.1 dated May 22, 2003). AMEC utilized the data usability criteria as defined in "MCP Representativeness Evaluations and Data Usability Assessments" (BWSC Policy # WSC-07-350), as well as the individual analytical methods defined in MADEP's Compendium of Analytical Methods (CAM,) to assess data quality and data usability. Data are considered to meet the requirements for defensibility, precision, accuracy and reporting of data are of sufficient quality to support this RAO unless otherwise stated below.

Groundwater Analytical Data Usability

This data usability assessment includes groundwater samples collected by Weston Solutions, Inc. The groundwater samples were collected on June 5, 2007; June 25, 2007; July 23, 2007; December 11, 2007; and May 19, 2008. The analytical data review elements included, but were not limited to laboratory control samples (LCS), surrogate recoveries, matrix spike/matrix spike duplicate (MS/MSD) recoveries, blank results, laboratory duplicates, field duplicates and laboratory case narratives.

Summary of Analytical Data Qualifications

Sample	Parameter	Issue	Use Limitation
DD-MW-002-R01-X		Di-n-butyl phthalate and bis(2-	
DD-MW-204-R02-X		ethylhexyl)phthalate were detected in	AMEC U-qualified all the
DD-MW-203-R04-X	8270	the method blank below the reporting	detected values that were <
DD-MW-207-R04-X		limits at 1.2 μg/L and 0.77 μg/L,	5X the blank concentrations.
DD-MW-205-R03-X		respectively.	
DD-MW-002-R01-X			AMEC III avalified all acid
DD-MW-204-R02-X		The assume note in board of recovered	AMEC UJ-qualified all acid
DD-MW-203-R04-X	8270	The surrogate phenol-d5 recovered	extractable analytes in these
DD-MW-207-R04-X		low in all associated samples.	samples due to the potential
DD-MW-205-R03-X			low bias.
DD-MW-002-R01-X			
DD-MW-204-R02-X		Aniline and phenol recovered low in	AMEC UJ-qualified aniline
DD-MW-203-R04-X	8270	the LCS/LCSD at 35%/27% and	and phenol in all samples
DD-MW-207-R04-X		12%/12%, respectively.	due to the low bias.
DD-MW-205-R03-X			
		Di-n-butyl phthalate and bis(2-	
DD-MW-206-R04-X		ethylhexyl)phthalate were detected in	AMEC U-qualified all the
DD-MW-201-R04-X	8270	the method blank below the reporting	detected values that were <
DD-MW-001-R02-X		limits at 0.89 μg/L and 0.48 μg/L,	5X the blank concentrations.
		respectively.	
DD-MW-206-R04-X		Aniline and phenol recovered low in	AMEC UJ-qualified aniline
DD-MW-201-R04-X	8270	the LCS/LCSD at 36%/32% and	and phenol in all samples
DD-MW-001-R02-X		12%/12%, respectively.	due to the low bias.

Sample	Parameter	Issue	Use Limitation
DD-MW-201-R04-X	8270	Aniline and phenol recovered low in the MS/MSD at 33%/26% and 12%/12%, respectively.	AMEC UJ-qualified aniline and phenol in this sample due to the low bias.
DD-MW-206-R04-X DD-MW-201-R04-X DD-MW-001-R02-X	8270	The surrogate phenol-d5 recovered low in all associated samples.	AMEC UJ- or J-qualified all acid extractable analytes in these samples due to the potential low bias.
DD-MW-206-R04-X DD-MW-201-R04-X DD-MW-001-R02-X	Mercury	Mercury was detected in the method blank below the reporting limit at 0.12 µg/L.	AMEC U-qualified all the detected values that were < 5X the blank concentrations.
DD-MW-201-R04-X	Metals	Arsenic and selenium recovered high in the MS/MSD at 128%/137% and 138%/156%, respectively.	AMEC J-qualified both arsenic and selenium in this sample due to the potential high bias.
DD-MW-208-R01-001-X DD-MW-208-R01-001-D	8270	Di-n-butyl phthalate and bis(2- ethylhexyl)phthalate were detected in the method blank below the reporting limits at 0.37 μg/L and 0.14 μg/L, respectively.	AMEC U-qualified all the detected values that were < 5X the blank concentrations.
DD-MW-208-R01-001-X DD-MW-208-R01-001-D	8270	The surrogate phenol-d5 recovered <10% in both samples.	AMEC R-qualified all acid extractable analytes in these samples due to the potential low bias.
DD-MW-208-R01-001-X DD-MW-208-R01-001-D	8270	Phenol recovered low in the LCS/LCSD at 13%/15%. 2,4-Dimethylphenol and 2,4-dinitrophenol had elevated RPDs at 21% and 36%, respectively.	These analytes were previously R-qualified by AMEC and have not been further qualified.
DD-MW-208-R01-001-X DD-MW-208-R01-001-D	8260	2,2-Dichloropropane and carbon disulfide recovered low in the LCSD at 66% and 67%.	AMEC UJ-qualified both compounds in these samples due to the low bias.
DD-MW-208-R01-001-X DD-MW-208-R01-001-D	ЕРН	Sample DD-MW-208-R01-001-D was submitted as a field duplicate of sample DD-MW-208-R01-001-X. Phenanthrene, C11-C22 aromatic ranges and total EPH had elevated RPDs.	AMEC J-qualified these analytes in the primary sample and its field duplicate due to the imprecision.
DD-MW-208-R02-X	8270	Di-n-butyl phthalate and bis(2- ethylhexyl)phthalate were detected in the method blank below the reporting limits at 0.77 μg/L and 0.70 μg/L, respectively.	AMEC U-qualified all the detected values that were < 5X the blank concentrations.
DD-MW-208-R02-X	8270	Phenol recovered low in the LCS/LCSD at 12%/13%.	AMEC UJ-qualified phenol in this sample due to the low bias.

Sample	Parameter	Issue	Use Limitation
DD-MW-208-R02-X	8270	The surrogate phenol-d5 recovered low in this sample.	AMEC UJ-qualified all acid extractable analytes in this sample due to the potential low bias.
DD-MW-207-R05-X DD-MW-207-R05-D DD-MW-208-R03-X DD-MW-203-R05-X DD-MW-201-R05-X	8270	Di-n-butyl phthalate and bis(2- ethylhexyl)phthalate were detected in the method blank below the reporting limits at 0.80 μg/L and 0.61 μg/L, respectively.	AMEC U-qualified all the detected values that were < 5X the blank concentrations.
DD-MW-207-R05-X DD-MW-207-R05-D DD-MW-208-R03-X DD-MW-203-R05-X DD-MW-201-R05-X	8270	Phenol recovered low in the LCS/LCSD at 23%/25%.	AMEC UJ-qualified phenol in all samples due to the low bias.
DD-MW-207-R05-X DD-MW-207-R05-D DD-MW-208-R03-X DD-MW-203-R05-X DD-MW-201-R05-X	Antimony	Antimony was detected in the method blank below the reporting limit at 0.78 µg/L.	All associated samples are ND and not impacted by the high bias.
DD-MW-002-R02-X DD-MW-201-R06-X DD-MW-201-R06-D DD-MW-204-R05-X DD-MW-205-R05-X	8270	The surrogate phenol-d5 recovered low in all associated samples.	AMEC UJ- or J qualified all acid extractable analytes in these samples due to the potential low bias.
DD-MW-206-R05-X DD-MW-002-R02-X DD-MW-201-R06-X DD-MW-201-R06-D DD-MW-203-R06-X DD-MW-204-R05-X DD-MW-205-R05-X DD-MW-206-R05-X DD-MW-207-R06-X DD-MW-208-R04-X	8270	Phenol (24%/25%), 4-nitrophenol (LCS 26%), and 1,4-dioxane (18%/20%) recovered low in the LCS/LCSD. Also 2,4-dinitrophenol had an elevated RPD of 46%.	AMEC UJ-qualified phenol, 4-nitrophenol, and 1,4- dioxane in all associated samples due to the low recovery. AMEC J-qualified 2,4-dinitrophenol in sample DD-MW-204-R05-X. This analyte was ND in all other samples and not impacted.
DD-MW-203-R06-X	8270	Aniline (29%/37%), phenol (20%/20%), 4-nitrophenol (MSD 28%), and 1,4-dioxane (14%/16%) recovered low in the MS/MSD. Also 4-chloroaniline had an elevated RPD of 26%.	AMEC UJ-qualified aniline, phenol, 4-nitrophenol, and 1,4-dioxane in sample DD-MW-203-R06-X due to the low bias. 4-Chloroaniline was ND and not impacted by the non-directional bias.
DD-MW-201-R06-X DD-MW-201-R06-D	Chromium	Sample DD-MW-201-R06-D was submitted as a field duplicate of sample DD-MW-201-R06-X. Chromium had an elevated RPD of 50%.	AMEC J-qualified chromium in the primary sample and its field duplicate due to the imprecision.

Sample	Parameter	Issue	Use Limitation
DD-MW-002-R02-X DD-MW-201-R06-X DD-MW-201-R06-D DD-MW-203-R06-X DD-MW-204-R05-X DD-MW-205-R05-X DD-MW-206-R05-X DD-MW-207-R06-X DD-MW-208-R04-X	Antimony	Antimony was detected in the method blank below the reporting limit at 0.37 µg/L.	AMEC U-qualified the detected antimony value in sample DD-MW-207-R06-X since the concentration was < 5X the blank concentration.
DD-MW-203-R06-X	Antimony	Antimony recovered low in the MS at 70%.	AMEC UJ-qualified antimony due to the potential low bias.

Groundwater Field Data Usability

The field component of the Data Usability Assessment evaluates whether the sampling procedure ensures that the samples collected and delivered to the laboratory are representative of each sampling point. The review elements included, but were not limited to appropriate sample collection procedures, holding times, sample receipt, appropriate sample containers, and sample preservation. Weston collected groundwater samples at the Site for laboratory analysis. The samples were kept in coolers with ice and submitted to the laboratory. All samples were analyzed within the required holding times specific for the analysis.

Groundwater data are considered to be usable under the MCP. The data are scientifically valid and defensible, and of a sufficient level of precision, accuracy, and completeness to support this RAO, with the exception of the acid extractable analytes in samples DD-MW-208-R01-001-X and DD-MW-208-R01-001-D.