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# REPORT

March 2021

## Immediate Response Action Completion

RTN 3-36304  
618R Waverly Street  
Framingham, MA

## TABLE OF CONTENTS

	Page
TABLE OF CONTENTS .....	i
LIST OF FIGURES.....	ii
LIST OF TABLES.....	ii
LIST OF APPENDICES.....	ii
1.0 INTRODUCTION.....	1-1
1.1 Site Description .....	1-2
1.2 Oil and Hazardous Materials (OHM) Usage and Storage .....	1-2
1.3 Area Receptors .....	1-2
1.4 Release Description .....	1-2
1.5 Reason for the IRA.....	1-4
1.6 Person(s) Undertaking the IRA.....	1-5
1.7 Activities Completed under the IRA .....	1-5
1.8 Future MCP Response Actions .....	1-6
2.0 IRA COMPLETION.....	2-1
3.0 LIMITATIONS.....	3-1

.....

## LIST OF FIGURES

Figure 1 ..... Locus Map  
Figure 2 ..... Sampling Plan  
Figure 3 ..... Area Receptors Map

## LIST OF TABLES

Table 1 ..... Summary of Soil Analytical Results – April 2020 Soil Borings  
Table 2 ..... Summary of Soil Analytical Results – Shallow Lead Concentrations

## APPENDICES

Appendix A ..... Public Notification Letters  
Appendix B ..... Laboratory Analytical Reports  
Appendix C ..... Risk Characterization Tables  
Appendix D ..... Photographic Log  
Appendix E ..... IDW Disposal Documents

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## 1.0 INTRODUCTION

Weston & Sampson Engineers, Inc. (Weston & Sampson) has prepared this Immediate Response Action (IRA) Completion Report for Release Tracking Number (RTN) 3-36304. This IRA Completion Report has been prepared on behalf of the City of Framingham, Massachusetts (the City), the current owner of Parcels 134-64-7867 (618R Waverly Street) and 134-64-9905 (no address) in Framingham, Massachusetts, and is being submitted in accordance with section 310 CMR 40.0427 of the Massachusetts Contingency Plan (MCP). The applicable Bureau of Waste Site Cleanup (BWSC) form, BWSC-105, has been submitted electronically, and copies of the public notification letters are presented in Appendix A.

As detailed in Section 1.4 of this Report, RTN 3-36304 is associated with the detection of lead in shallow soil at concentrations indicative of an Imminent Hazard (IH) in an area comprising approximately 5,900 square feet straddling Parcels 134-64-7867 and 134-64-9905 (the Site); this finding is referred to as a "release," as described by the MCP at 310 CMR 40.0950. The Massachusetts Department of Environmental Protection (MassDEP) was verbally notified of the 2-hour release condition on June 3, 2020. MassDEP subsequently issued RTN 3-36304 to the Site and approved an IRA including temporarily fencing the area around the two locations that constituted the IH condition with a 6-foot high fence, and placing warning signs on the fence. MassDEP also noted in additional comments at the end of the Release Log Form (Section H.16) that exposed soil should be covered with high-density polyethylene sheeting to prevent airborne dust; however, due to heavy vegetation (including trees, brush, and weeds/grasses) at the Site, installation of polyethylene sheeting on the ground surface to prevent airborne dust was not feasible or necessary. This was discussed with MassDEP (Allison Williams) at the time of Notification on June 3, 2020.

An IRA Plan for RTN 3-36304 was submitted to the MassDEP on July 31, 2020. Installation of approximately 320 linear feet of 6-foot-tall, temporary chain-link fencing around the two locations that constituted the IH was completed on June 15, 2020 and signs were posted in three languages (English, Spanish, and Portuguese) warning of lead in the soil as a potential hazard inside the fenced area. Based on the sampling data, the IH area is estimated to encompass approximately 5,900 square feet around two sample locations with elevated lead concentrations. Temporary fencing was installed to prevent access to the IH area until further assessment activities are completed.

On September 11, 2020, the City submitted a Release Notification Form (RNF) for the detection of metals, polycyclic aromatic hydrocarbons (PAHs), and volatile organic compounds (VOCs) in soil at concentrations above the Reporting Category S-1 (RCS-1) on all of Parcels 134-64-7867 (618R Waverly Street) and 134-64-9905 (no address). MassDEP assigned RTN 3-36463 to the 120-day reporting condition. The Disposal Site Boundary (DSB) for RTN 3-36464 includes the DSB for RTN 3-36304.

In emails dated November 19 and December 1, 2020, MassDEP confirmed to the City their position that it is appropriate to submit an IRA Completion Report for RTN 3-36304. MassDEP and the City agree that additional sampling is required to evaluate the extent of lead in shallow soil at the Site. This sampling will be completed as part of future Comprehensive Response Actions at the Site subsequent to the submittal of this IRA Completion Report. It is the intent of the City to link the two RTNs (3-36304 and 3-36463) through Tier Classification, and perform Comprehensive Response Actions for the Site under RTN 3-36463.

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## 1.1 Site Description

The Site consists of an approximately 5,900 square foot area on two parcels: a larger parcel with Tax ID 134-64-7867 (currently known and numbered as 618R Waverly Street) and a smaller parcel with Tax ID 134-64-9905 (no address) located in Framingham, Massachusetts. The City owns both parcels, which together cover approximately 2.04 acres of vacant land comprised primarily of woodlands and wetland, with a paved area to the north and a paved area to the west.

The property at 618R Waverly Street is a vacant property owned by the City comprising approximately 1.66 acres. The adjacent approximately 0.38-acre parcel to the east is also owned by the City and identified as Parcel 134-64-9905; the Site covers a portion of each parcel. Figure 1 shows the Site and general surroundings. Figure 2 depicts the Site in more detail, and includes soil boring locations, property boundaries, and other pertinent information.

The Site is primarily woodlands, with areas of wetland abutting to the south and parking lot abutting to the north. A small path-like area that local residents may use to enter areas of Cedar Woods bisects the Site north-to-south. During response actions described below, personal belongings of a trespasser were observed outside of the IH area, indicating likely use/occupation of the property by at least one person experiencing housing insecurity.

Waverly Street and commercial businesses are located north of the Site, to the east is undeveloped land with mixed commercial and residential properties along Waverly Court, to the south are wetlands and wooded parcels owned by the City (generally referred to as "Cedar Woods"), and to the west is Mellen Street with commercial properties beyond the street to the west. There are no surface water bodies in the immediate vicinity of the Site; Farm Pond is located approximately 600 feet north of the Site and a wetland area 50-100 feet south of the Site. The Site boundary is not fenced and is accessible to the public as open space.

## 1.2 Oil and Hazardous Materials (OHM) Usage and Storage

Weston & Sampson has not observed any evidence of the existence, past or present, of underground storage tanks (USTs). No other current or historical OHMs storage has been identified at the Site. The source of the elevated lead levels is unknown but is likely associated with fill material and/or past uses of the property and surrounding properties.

## 1.3 Area Receptors

As presented in Figure 3, Area Receptors Map, the Site is not located within a Zone II or Interim Well Protection Area (IWPA) for a public water supply. The Site is also not located in a potentially productive aquifer or within a Current or Potential Drinking Water Source Area. In addition, there are no known drinking water wells within 500 feet of the Site. No portion of the Site is listed as Natural Heritage and Endangered Species Protected (NHESP) Wetlands Habitats for Rare or Endangered Species. The majority of the Site, except for the paved western portion of the Site and the eastern half of the Waverly Street – paper road parcel, is identified as an area of Protected Open Space. The closest surface water bodies to the Site are Farm Pond, which is located 600 feet north of the Site and a wetland area, which is 50 to 100 feet south of the Site.

## 1.4 Release Description

On April 6, 7 and 8, 2020, Weston & Sampson oversaw the advancement of 15 soil borings (SB-101 through SB-115) on the two above-mentioned City-owned parcels. The work was completed as part of a US EPA Brownfields-funded environmental assessment. The soil borings were advanced by New

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England Geotech of Jamestown, Rhode Island (NE Geotech) using a mix of direct-push technology and hand tooling. Soil samples were collected from each soil boring at the 0-1 foot (ft) and 0-3 ft depths below grade. Additional samples were collected at varying depths, depending on field evidence of impacts, to a maximum depth of 14 ft below grade. Soil samples were transported via courier under chain-of-custody to Absolute Resource Associates of Portsmouth, New Hampshire (ARA). Soil analytical results are summarized in Tables 1 and 2, and ARA laboratory analytical reports are included as Appendix B.

An initial set of soil samples, including 0-3 ft samples from all locations, were analyzed by ARA for Massachusetts Compendium of Analytical Methods (CAM)-14 Metals (antimony, arsenic, barium, beryllium, cadmium, chromium (III), chromium (VI), lead, mercury, nickel, silver, thallium, vanadium, zinc), extractable petroleum hydrocarbons (EPH) with target PAHs, volatile petroleum hydrocarbons (VPH) with target VOCs, and/or VOCs by state or US EPA methods. Soil samples collected for metals analysis from the 0-1 ft depth interval were held by the laboratory pending the receipt and review of the initial dataset of soil from the 0-3 ft depth interval.

Laboratory analytical results for the initial dataset identified lead at concentrations in excess of the MCP Method 1 S-1 Cleanup Standard of 200 mg/kg in 5 of the 0-3 foot samples collected from unpaved locations (SB-105, SB-107, SB-108, SB-109 and SB-111). The Method 1 S-1 Cleanup Standard was used as a conservative action level to evaluate whether activation of the remaining shallow soil samples collected from unpaved locations was necessary. Concentrations of lead in the 0 to 3 feet sampling intervals ranged from 630 to 1,300 mg/kg. Based on these results, it was unclear if a potential IH condition was present; therefore, on May 15, 2020, Weston & Sampson instructed ARA to activate the shallow (0-1 ft) soil samples.

Eight (8), 0-1 foot samples from the locations where access to the soil is not prevented by means of a physical barrier, such as pavement (SB-102, SB-106, SB-107, SB-108, SB-109, SB-110, SB-111, and SB-112), were activated for laboratory analysis of lead. The resulting concentrations of lead in soil from the 0-1 foot interval ranged from 11 to 900 mg/kg. Two samples, SB-108 (0-1) and SB-109 (0-1), from the wooded area in approximately the center of the two parcels, had concentrations of 900 and 790 mg/kg, respectively, which were an order of magnitude greater than lead in the surrounding samples, ranging from 11 to 73 mg/kg. Due to the presence of lead in exposed surface soil (0-1 foot) at concentrations approximately an order of magnitude above the surrounding samples, SB-108 and SB-109 were considered to be a "Hot Spot," as defined by the MCP, and were further considered in an IH Evaluation as described below.

#### IH Evaluation

Weston & Sampson received the laboratory report and reviewed the results for the eight (8) 0-1' soil samples on June 2, 2020, and performed an Imminent Hazard Evaluation for the Site.

#### *Potential Human Receptors*

Potential human receptors at the Site include the following:

- Park Visitors - Adults and children may be exposed to impacted Site soil via incidental ingestion and dermal exposure.
- Utility Workers - Adults conducting potential repairs/maintenance activities may be exposed to impacted Site soil via incidental ingestion, dermal contact, and inhalation of fugitive dust.

- Construction - Adults conducting potential future subsurface construction activities may be exposed to impacted Site soil via incidental ingestion, dermal contact, and inhalation of fugitive dust.

MassDEP's spreadsheets (ShortForms) for assessing potential risks to human health were used to evaluate the data for the lead results. These ShortForms use various exposure scenarios to evaluate whether or not the Site conditions meet the MCP definition of a potentially significant risk to human health. Weston & Sampson discussed with the City two ShortForm scenarios for the IH evaluation at the Site, either using the "Park Visitor" exposure (younger children present at the property) or the "Trespasser" (older children and adults) exposure scenario. The City, in discussion with Weston & Sampson, recommended evaluating the potential exposure at the Site using the Park Visitor scenario since the Site is mapped as "open space;" in addition, there is a small path-like area that local residents that may use as way to access areas of Cedar Woods, evidence of persons experiencing housing insecurity at this location, and the potential for unrestricted use by nearby residents or visitors. This approach was very conservative to protect potential public health exposures, even though the Site does not currently contain any park amenities and is not considered an active park.

#### *Exposure Point Concentrations (EPC)*

For potential park visitor (the most sensitive of the receptors identified) exposures to lead in the two locations sampled at the Site, the average concentration of SB-108 and SB-109, 845 mg/kg, was used to represent potential exposures.

#### *Risk Calculations*

Local residents that may use this portion of Cedar Woods for recreation or as way to enter other areas of Cedar Woods were assumed to be exposed to the lead EPC via dermal contact and incidental ingestion of soil. Conservatively, the MassDEP Shortform (October 2012 with updates March 2015) for park visitor exposure to soil was used to evaluate local residential exposures to soil and included standard exposure assumptions. The risk characterization tables were included with the IRA Plan.

The estimated risks for the park visitor age 1 to 2 years (noncancer, imminent hazard exposure) were greater than MassDEP risk limit of 1 (2.4). Since the Imminent Hazard Limit for lead is also 1, the results of this Imminent Hazard evaluation indicated that there is a potential Imminent Hazard associated with exposure to soil represented by SB-108 and SB-109. Risk characterization tables are included in Appendix C.

### **1.5 Reason for the IRA**

In accordance with 310 CMR 40.0412(1) through (4), an IRA is required at the Site because a release of OHM occurred which required notification to MassDEP under both the "2 hour" notification provision of the Massachusetts Contingency Plan (MCP) and because the release was assessed to pose an IH. Pursuant to MCP Section 40.0412, IRAs shall be conducted at the following sites:

- (1) Sites or vessels where a release or threat of release of oil and/or hazardous material has occurred which requires notification to the Department under the "Two-Hour" notification provisions of 310 CMR 40.0311 or 40.0312;
- (2) Sites where a release or threat of release of oil and/or hazardous material has occurred which requires notification to the Department under the "72-Hour" notification provisions of 310 CMR 40.0313 or 40.0314;

- (3) Sites where a release of oil and/or hazardous material has resulted in conditions which have been determined to pose an Imminent Hazard pursuant to 310 CMR 40.0950; and
- (4) Any other site or vessel where the Department determines that immediate or accelerated response actions are necessary to prevent, eliminate, or minimize damage to health, safety, public welfare or the environment.

As described in detail above, the results of an IH Evaluation indicated that Site conditions (i.e., lead concentrations in shallow soil) did pose an IH because access to the lead-impacted Site soil was unrestricted.

In accordance with the MCP requirements for release notification, the MassDEP was initially notified of the IH condition on June 3, 2020, within 2 hours of the owner first having knowledge of the site conditions. The MassDEP subsequently assigned RTN 3-36304 to this reportable condition, and verbally approved an IRA including temporarily fencing the area around the two locations that constituted the IH condition (SB-108 and SB-109) with a 6-foot high fence and placing warning signs on the fence.

In accordance with the MCP requirements specified at 310 CMR 40.0426(2), an IH Evaluation was initiated within fourteen days of obtaining knowledge of such a condition. As previously mentioned, the risk calculations utilizing MassDEP Short Forms for park visitors confirmed that an IH condition existed at the Site.

### 1.6 Person(s) Undertaking the IRA

The party assuming responsibility for conducting the IRA, and the current owner of the Site is:

Dr. Yvonne M. Spicer, Mayor  
City of Framingham  
150 Concord Street  
Framingham, Massachusetts 01702  
mayor@framinghamma.gov  
(508) 532-5400

### 1.7 Activities Completed under the IRA

#### Installation and Repair of Temporary Fence

On June 15, 2020, Weston & Sampson observed the installation of approximately 320 linear feet of 6-foot-tall, temporary chain-link fencing in the area around SB-108 and SB-109, as shown in Figure 2. Installation of the fence was completed by East Coast Fence of Kingston, Massachusetts (ECF) under contract with the City. The fencing was installed as a short-term risk mitigation measure to restrict access to the surficial soil represented by shallow samples collected at SB-108 and SB-109. Signs were posted in three languages (English, Spanish, Portuguese) as warning of lead in the soil as a potential hazard inside the fenced area. Photographs of the completed fencing installation and signage are included as Appendix D. Due to heavy vegetation (including trees, brush, and weeds/grasses) at the Site, it was determined that installation of polyethylene sheeting on the ground surface to prevent airborne dust was not feasible or necessary. This was discussed with MassDEP (Allison Williams) at the time of Notification on June 3, 2020.



MassDEP personnel (Allison Williams) of the ER group conducted a site visit on June 18, 2020. MassDEP identified two areas of fencing in need of repair. On June 19, 2020, Brian Smith of the Framingham Department of Public Health repaired the two holes by weaving wire through the fencing. The two damaged panels of fencing at the 618R Waverly Street IH area were then replaced on August 6, 2020.

Based on the sampling data, the IH area is estimated to encompass approximately 5,900 square feet around two sample locations with elevated lead concentrations. Temporary fencing was installed to prevent access to the IH area.

#### Management of Remediation Waste

Investigation Derived Waste (IDW), generated during the installation of the soil borings in April 2020, was managed as Remediation Waste due to the presence of RCS-1 exceedances as described above in section 1.4. During the course of drilling activities, approximately 0.2 cubic yards of surplus drill cuttings were placed in a labeled and secured 55-gallon steel drum and stored on-site. The IDW was generated in small volumes from each of the boring locations depicted on the attached Figure 2; as such, it is represented by the laboratory analytical results described above.

Subsequent to the April subsurface investigation, on June 15, 2020, Weston & Sampson collected one composite sample from the drummed soil (IDW-1) and submitted it to ARA for analysis of arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver by Toxicity Characteristic Leaching Procedures (TCLP) methods. Copies of the laboratory analytical reports were previously included with the IRA Plan.

On July 29, 2020, the drummed IDW was transported under a Bill of Lading by Strategic Environmental Services of Sutton, Massachusetts (SES) to Northland Environmental in Providence, Rhode Island (Northland), an appropriately permitted disposal facility in Rhode Island. This facility will temporarily store the IDW, combine it with similar material, and dispose of it at a licensed landfill, which is an appropriate disposal option based on the representative analytical data. Supporting documentation for the transportation and disposal of the IDW is included in Appendix E.

### **1.8 Future MCP Response Actions**

#### Maintenance and Monitoring

City personnel continue to conduct periodic inspections of the temporary fencing installed to mitigate the IH condition at the Site. Should fencing require repair or replacement, the City will contact the fencing contractor to perform maintenance as required. The fencing will remain in place until an updated IH Evaluation indicates that an IH condition does not exist without the fence in place.

#### Assessment

To further evaluate the extent of lead in shallow soil in the area surrounding SB-108 and SB-109, additional soil samples will be collected as part of future Comprehensive Response Actions conducted at the Site.

#### Remedial Activities

The results of additional soil sampling in the area surrounding SB-108 and SB-109 will inform future remedial activities. Future remedial activities will be conducted as Comprehensive Response Actions in accordance with the MCP.

## 2.0 IRA COMPLETION

The MCP requirements for an IRA Completion Report, pursuant to 310 CMR 40.0427, are presented below in ***bold, italic text***, with the corresponding required information provided in normal text.

***(a) a description of the release or threat of release, site conditions and surrounding receptors;***

Descriptions of the Release, Site conditions, and surrounding receptors are included in Sections 1.4, 1.1-1.2, and 1.3, respectively.

***(b) a description of the work completed, including work undertaken in response to any conditions of approval imposed by the Department, and any work undertaken at the site that was not included in the scope of the Immediate Response Action Plan, where submitted;***

A description of the work completed under the IRA is included in Section 1.7.

***(c) all investigatory and monitoring data obtained during the implementation of the Immediate Response Action;***

No investigatory or monitoring data was collected during implementation of the IRA. Investigatory data collected prior to the implementation of the IRA is summarized in Section 1.4.

***(d) a succinct statement of the findings and conclusions of the Immediate Response Action;***

Based on the results of the 2020 subsurface investigation noted in Section 1.4 above and described in more detail in Sections 3 and 4, Weston & Sampson performed an IH Evaluation for a lead “Hot Spot” in shallow soil at the Site. The results of the IH Evaluation indicated an IH condition was present at the Site. The City of Framingham notified MassDEP of the IH condition on June 3, 2020 within two hours of obtaining knowledge of the IH, and RTN 3-36304 was assigned to the Release.

Activity conducted under IRA, verbally approved by MassDEP at the time of notification and further documented in the July 2020 IRA Plan, have included the installation of approximately 320 linear feet of 6-foot tall chain-link fencing around the affected area and appropriate signage. As a result of the IRA activities conducted under RTN 3-36304, the IH condition was eliminated. The chain-link fence will remain in place until an updated IH Evaluation indicates that an IH condition does not exist without the fence in place.

***(e) details and documentation on the management of any Remediation Waste, Remedial Wastewater and/or Remedial Additives managed at the site as part of the Immediate Response Action;***

A summary of the management and off-site disposal of Remediation Waste is included in Section 1.7. Shipping documents are provided in Appendix E.

***(f) a description of any ongoing activities related to the Immediate Response Action that will be conducted at the site, including monitoring activities, security measures and the maintenance of fences, caps and other passive systems; and***

A summary of future activities at the Site related to the IRA, including the maintenance of installed fencing, is included in Section 1.8.

*(g) a description of any ongoing activities related to the Immediate Response Action that will be conducted at the site as part of Comprehensive Response Actions*

A summary of future activities at the Site related to the IRA, including additional sampling and potential remedial activities conducted as part of Comprehensive Response Actions, is included in Section 1.8.

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### 3.0 LIMITATIONS

This report was prepared for the use by the City of Framingham exclusively. The findings provided by Weston & Sampson in this report are based solely on the information reported in this document and previous documents prepared for the Site. Future investigations, and/or information that was not available to Weston & Sampson at the time of the investigation, may result in a modification of the findings stated in this report.

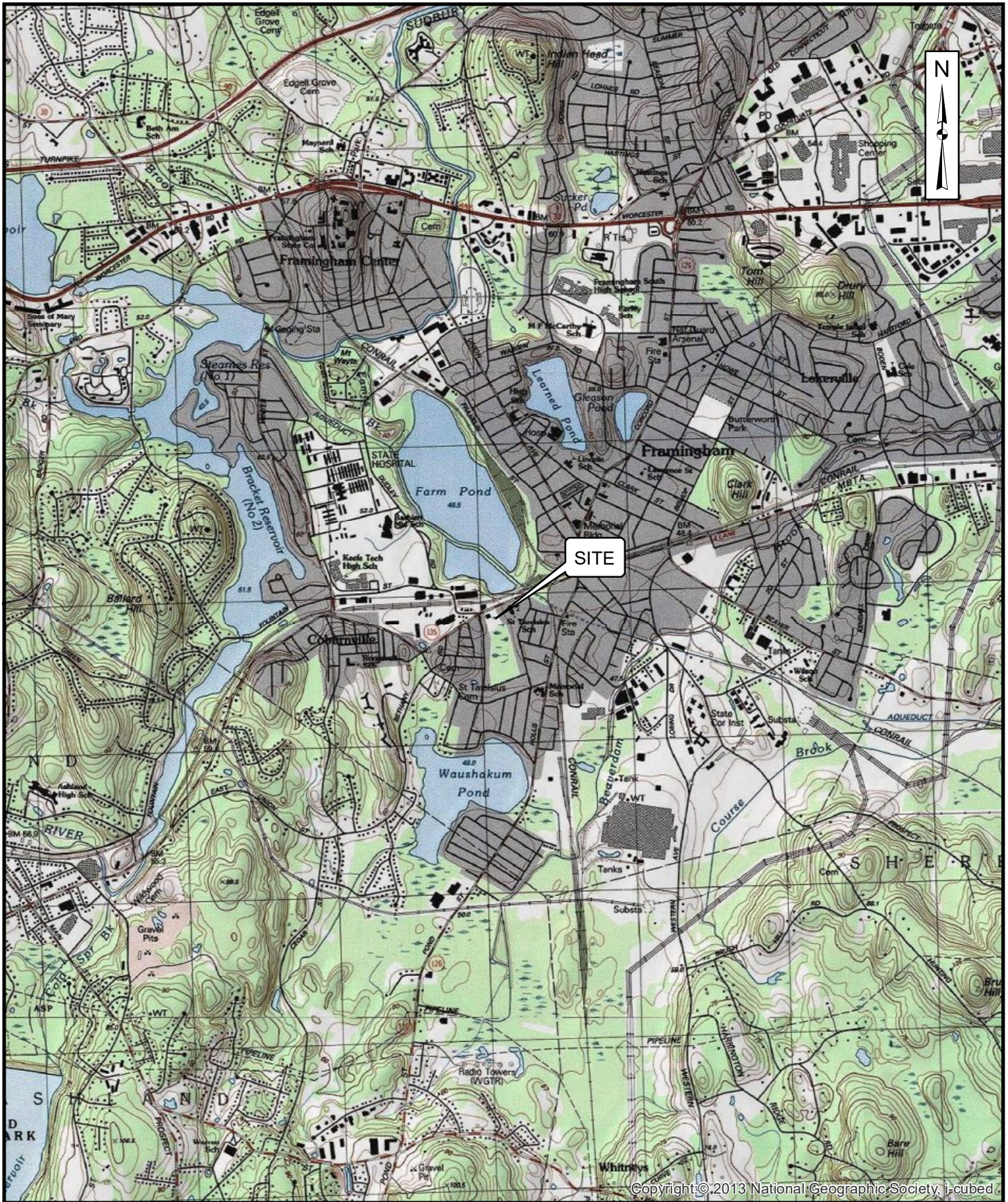
Additional information that becomes available concerning this Site or neighboring properties, which could directly impact the Site in the future, should be made available to Weston & Sampson for review so that, if necessary, conclusions presented in this report may be modified. The conclusions of this report are based on Site conditions observed by Weston & Sampson personnel at the time of the investigation, information provided by the City of Framingham, and samples collected and analyzed on the dates shown or stated in this report. This report has been prepared in accordance with generally accepted engineering and geological practices. No other warranty, express or implied, is made.

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JRS/FMR

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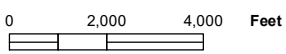
## FIGURES

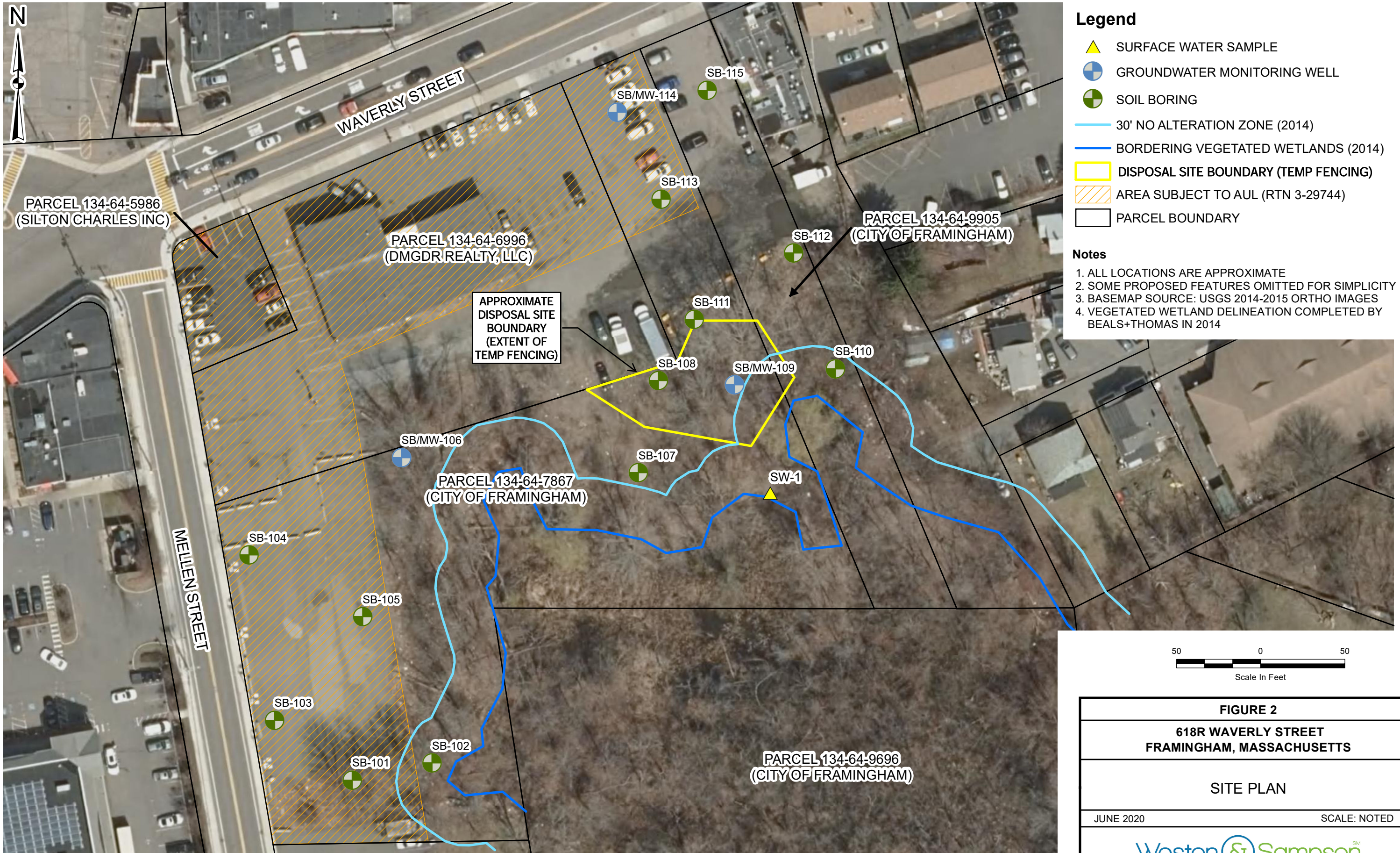
C:\Users\spencerj\OneDrive - WESTON & SAMPSON ENGINEERS, Inc\Documents\ArcGIS\Framingham\CW - locus.mxd -> 7/27/2020 6:02:27 PM -> 7/27/2020 6:03:16 PM











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**FIGURE 1**  
**618R WAVERLY STREET**  
**FRAMINGHAM, MASSACHUSETTS**  
**LOCUS MAP**



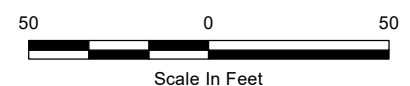



**Legend**

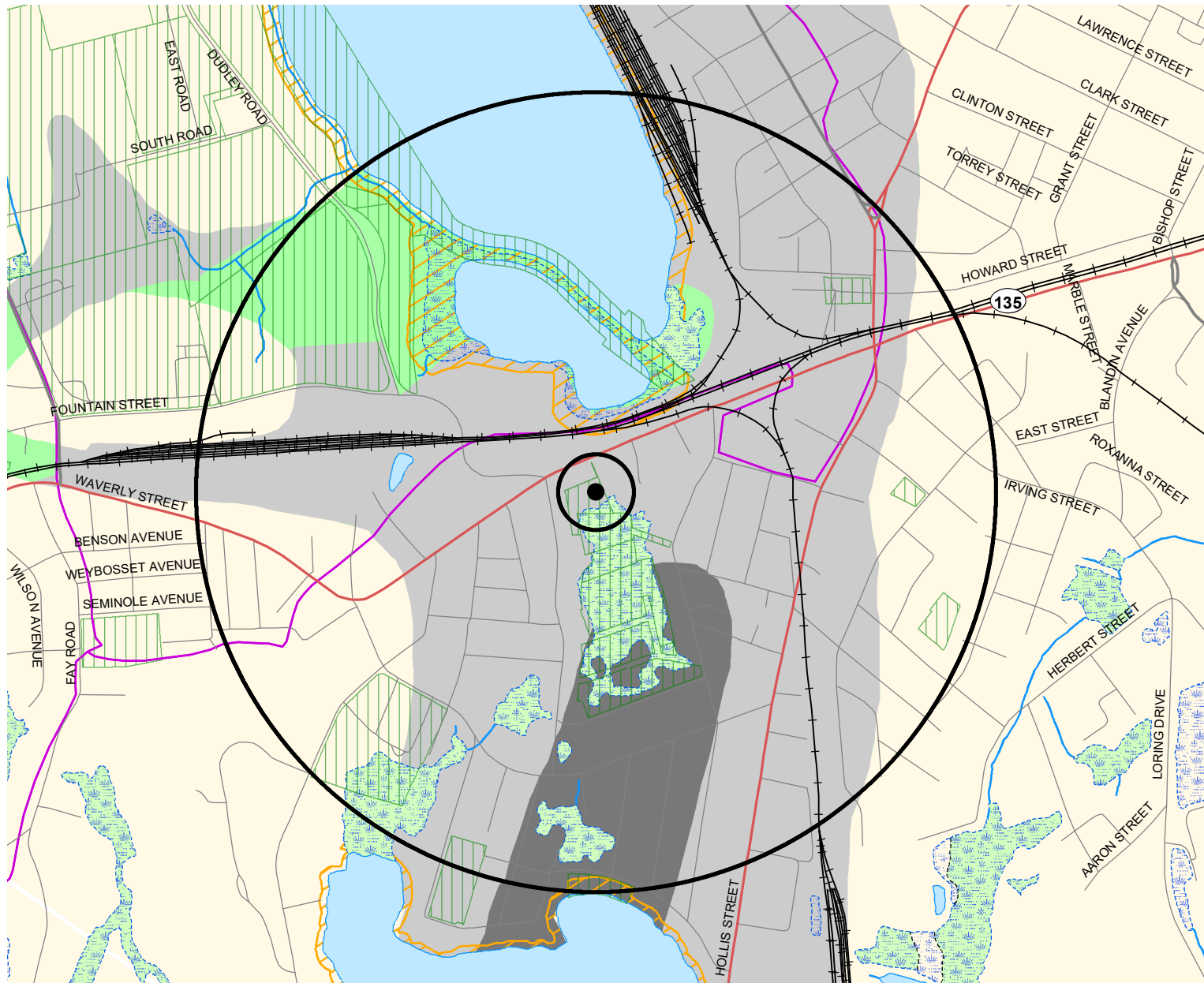
-  SURFACE WATER SAMPLE
-  GROUNDWATER MONITORING WELL
-  SOIL BORING
-  30' NO ALTERATION ZONE (2014)
-  BORDERING VEGETATED WETLANDS (2014)
-  DISPOSAL SITE BOUNDARY (TEMP FENCING)
-  AREA SUBJECT TO AUL (RTN 3-29744)
-  PARCEL BOUNDARY

**Notes**

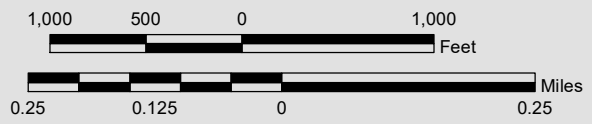
1. ALL LOCATIONS ARE APPROXIMATE
2. SOME PROPOSED FEATURES OMITTED FOR SIMPLICITY
3. BASEMAP SOURCE: USGS 2014-2015 ORTHO IMAGES
4. VEGETATED WETLAND DELINEATION COMPLETED BY BEALS+THOMAS IN 2014



<b>FIGURE 2</b>	
<b>618R WAVERLY STREET FRAMINGHAM, MASSACHUSETTS</b>	
SITE PLAN	
JUNE 2020	SCALE: NOTED
	



- Legend**
- ▲ Ground Water
  - ▲ Surface Water
  - ▲ Non-Community
  - ✱ NHESP Certified Vernal Pools
  - Railroads by Ownership
  - Pipeline
  - Pipeline Arbitrary Extension
  - Powerline
  - Powerline Arbitrary Extension
  - Ski Lift/Tramway
  - Substation
  - Landing Strip/Airport
  - ◆ Highway Exit Locations
- All Roads**
- Road Classification**
- Limited Access Highway
  - Multi-lane Hwy, not limited access
  - Other Numbered Highway
  - Major Road, Collector
  - Minor Road, Arterial
  - Sub-basins
  - Major Basins
  - Landfills
  - Dumping Grounds
  - Protected Open Space
  - ACECs
  - Zone A
  - IWPA
  - DEP Approved Zone IIs
  - Sole Source Aquifers
  - NHESP Estimated Habitats of Rare Wildlife
  - NHESP Priority Habitats of Rare Species
- Non Potential Drinking Water Source Area**
- High Yield
  - Medium Yield
- Aquifers**
- High Yield
  - Medium Yield
- MA Towns (from Survey Points)**
- MA Towns (from Survey Points)
  - Marsh/Bog
  - Wooded marsh
  - Cranberry Bog
  - Salt Marsh
  - Open Water
  - Reservoir (with PWSID)
  - Tidal Flats
  - Beach/Dune



**Data Source:** Office of Geographic and Environmental Information (MassGIS), Commonwealth of Massachusetts Executive Office of Environmental Affairs

**NOTE:** Radii shown are approximately 500 feet and 1/2 mile from Site boundary.

**FIGURE 3**  
 AREA RECEPTORS MAP  
 618R WAVERLY STREET  
 CITY OF FRAMINGHAM, MASSACHUSETTS



## TABLES

**Table 1**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**618R Waverly Street**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)		Sample Location, Depth, and Date															
					RCS-1		S-1/GW-2		S-1/GW-3		SB-101		SB-102		SB-103		SB-104		SB-105	
					0-3 feet	9-12 feet	0-3 feet	11-14 feet	0-3 feet	3-5 feet	7-10 feet	0-3 Feet	8-10 Feet	0-3 feet	5-8 feet					
					4/6/2020	4/6/2020	4/6/2020	4/6/2020	4/6/2020	4/6/2020	4/6/2020	4/7/2020	4/7/2020	4/6/2020	4/6/2020					
<b>EPH</b>																				
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<21	<25	<20	<46	NT	NT	NT	<20	<22	<24	<22					
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>35</b>	<b>82</b>	<b>130</b>	<b>80</b>	NT	NT	NT	<b>34</b>	<b>100</b>	<b>280</b>	<b>23</b>					
C11-C22 Aromatics	mg/kg	1000	1000	1000	<21	<b>41</b>	<b>50</b>	<46	NT	NT	NT	<b>25</b>	<22	<b>130</b>	<22					
<b>Target PAHS</b>																				
Acenaphthene	mg/kg	4	1000	1000	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Acenaphthylene	mg/kg	1	600	10	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Anthracene	mg/kg	1000	1000	1000	<0.21	<b>0.40</b>	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Benzo(A)Anthracene	mg/kg	7	7	7	<0.21	<b>0.98</b>	<0.20	<0.46	NT	NT	NT	<b>0.40</b>	<0.22	<b>1.5</b>	<0.22					
Benzo(A)Pyrene	mg/kg	2	2	2	<0.21	<b>0.89</b>	<0.20	<0.46	NT	NT	NT	<b>0.44</b>	<0.22	<b>1.4</b>	<0.22					
Benzo(B)Fluoranthene	mg/kg	7	7	7	<0.21	<b>0.88</b>	<0.20	<0.46	NT	NT	NT	<b>0.46</b>	<0.22	<b>1.6</b>	<0.22					
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<0.21	<b>0.72</b>	<0.20	<0.46	NT	NT	NT	<b>0.35</b>	<0.22	<b>1.2</b>	<0.22					
Benzo(K)Fluoranthene	mg/kg	70	70	70	<0.21	<b>0.71</b>	<0.20	<0.46	NT	NT	NT	<b>0.44</b>	<0.22	<b>1.2</b>	<0.22					
Chrysene	mg/kg	70	70	70	<0.21	<b>1.0</b>	<0.20	<0.46	NT	NT	NT	<b>0.56</b>	<0.22	<b>1.9</b>	<0.22					
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<b>0.38</b>	<0.22					
Fluoranthene	mg/kg	1000	1000	1000	<b>0.22</b>	<b>2.1</b>	<0.20	<b>0.48</b>	NT	NT	NT	<b>0.80</b>	<0.22	<b>3.2</b>	<0.22					
Fluorene	mg/kg	1000	1000	1000	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<0.21	<b>0.63</b>	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<b>1.1</b>	<0.22					
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Naphthalene	mg/kg	4	20	500	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22					
Phenanthrene	mg/kg	10	500	500	<0.21	<b>2.1</b>	<0.20	<0.46	NT	NT	NT	<b>0.39</b>	<0.22	<b>1.5</b>	<0.22					
Pyrene	mg/kg	1000	1000	1000	<b>0.24</b>	<b>2.3</b>	<0.20	<b>0.56</b>	NT	NT	NT	<b>0.85</b>	<0.22	<b>2.9</b>	<0.22					
<b>VPH</b>																				
C5-C8 Aliphatics	mg/kg	100	100	100	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3					
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3					
C9-C10 Aromatics	mg/kg	100	100	100	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3					
<b>Target VOCs</b>																				
Benzene	mg/kg	2	40	40	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
Ethylbenzene	mg/kg	40	500	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
Naphthalene	mg/kg	4	20	500	NT	<0.31	NT	<0.78	NT	NT	<0.29	NT	<0.24	NT	<0.22					
Toluene	mg/kg	30	500	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
m-p Xylene	mg/kg	100	100	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
o-Xylene	mg/kg	100	100	500	NT	<0.13	NT	<.31	NT	NT	<0.12	NT	<0.097	NT	<0.087					
<b>Metals</b>																				
Antimony	mg/kg	20	20	20	<0.52	<b>2.7</b>	<0.54	<1.1	<0.51	<b>3.7</b>	<0.60	<0.54	<0.57	<b>4.7</b>	<0.55					
Arsenic	mg/kg	20	20	20	<b>3.3</b>	<b>28</b>	<b>3.3</b>	<b>10</b>	<b>3.5</b>	<b>13</b>	<b>3.3</b>	<2.7	<b>3.1</b>	<b>7.3</b>	<b>3.0</b>					
Barium	mg/kg	1000	1000	1000	<b>18</b>	<b>140</b>	<b>27</b>	<b>56</b>	<b>25</b>	<b>250</b>	<b>36</b>	<b>17</b>	<b>27</b>	<b>310</b>	<b>23</b>					
Beryllium	mg/kg	90	90	90	<0.52	<0.68	<0.54	<1.1	<0.51	<0.62	<0.60	<0.54	<0.57	<0.60	<0.55					
Cadmium	mg/kg	70	70	70	<0.52	<b>1.3</b>	<0.54	<1.1	<0.51	<b>4.9</b>	<0.60	<0.54	<0.57	<b>2.0</b>	<0.55					
Chromium (III)	mg/kg	1000	1000	1000	NT	<b>51</b>	NT	<b>41</b>	NT	<b>9.4</b>	NT	NT	<b>16</b>	<b>18</b>	NT					
Chromium (VI)	mg/kg	100	100	100	<b>7.3</b>	<b>3.8</b>	<b>17</b>	<1.00	<b>6.9</b>	<0.49	<b>18</b>	<5.4	<0.47	<0.48	<b>11</b>					
Lead	mg/kg	200	200	200	<b>8.7</b>	<b>230</b>	<b>16</b>	<b>190</b>	<b>22</b>	<b>710</b>	<b>8.4</b>	<b>16</b>	<b>5.5</b>	<b>650</b>	<b>5.0</b>					
Mercury	mg/kg	20	20	20	<0.15	<b>0.58</b>	<0.13	<0.31	<0.15	<b>0.58</b>	<0.17	<0.15	<0.16	<b>2.9</b>	<0.15					
Nickel	mg/kg	600	600	600	<b>7.4</b>	<b>14</b>	<b>23</b>	<b>64</b>	<b>6.3</b>	<b>12</b>	<b>13</b>	<b>5.4</b>	<b>17</b>	<b>18</b>	<b>14</b>					
Selenium	mg/kg	400	400	400	<5.2	<6.8	<5.4	<11	<5.1	<6.2	<6.0	<5.4	<5.7	<6.0	<5.5					
Silver	mg/kg	100	100	100	<2.6	<3.4	<2.7	<5.6	<2.5	<3.1	<3.0	<2.7	<2.8	<3.0	<2.7					
Thallium	mg/kg	8	8	8	<0.52	<0.68	<0.54	<1.1	<0.51	<0.62	<0.60	<0.54	<0.57	<0.60	<0.55					
Vanadium	mg/kg	400	400	400	<b>11</b>	<b>15</b>	<b>20</b>	<b>17</b>	<b>10</b>	<b>10.0</b>	<b>25</b>	<b>8.3</b>	<b>22</b>	<b>19</b>	<b>14</b>					
Zinc	mg/kg	1000	1000	1000	<b>17</b>	<b>220</b>	<b>31</b>	<b>530</b>	<b>19</b>	<b>920</b>	<b>18</b>	<b>16</b>	<b>18</b>	<b>770</b>	<b>69</b>					
<b>VOCs</b>																				
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Cis-1,2-Dichloroethylene	mg/kg	0.1	0.1	100	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT					

QC by JMR 5/8/2020

**Abbreviations:**

EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
PAH = Polycyclic Aromatic Hydrocarbons  
NT = Not Tested  
BDL = Below Detection Limit  
MCP = Massachusetts Contingency Plan  
mg/kg = milligram per kilogram

**Notes:**

-- = No standard available  
< = indicates parameter not detected above laboratory method reporting limit, shown  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter equals or exceeds the applicable MCP RCS-1  
**BOLD** Parameter exceeds the applicable MCP RCS-1 and the Method 1 S-1 Cleanup Standard  
1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.

**Table 1**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**618R Waverly Street**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)		Sample Location, Depth, and Date														
					RCS-1	S-1/GW-2	S-1/GW-3	SB-106			SB-107		SB-108		SB-109		SB-110		
								0-3 feet	3-4 feet	5-8 feet	0-3 Feet	10-11 Feet	0-3 Feet	6-9 Feet	0-3 Feet	5-8 Feet	DUP-1	0-3 Feet	5-8 Feet
								4/6/2020	4/6/2020	4/6/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020
<b>EPH</b>																			
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<21	NT	<240	<23	<b>34</b>	<26	<b>92</b>	<28	<78	<83	<22	<110			
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>91</b>	NT	<b>570</b>	<b>36</b>	<b>410</b>	<b>39</b>	<b>1900</b>	<b>42</b>	<b>290</b>	<b>390</b>	<b>68</b>	<b>140</b>			
C11-C22 Aromatics	mg/kg	1000	1000	1000	<b>44</b>	NT	<b>290</b>	<23	<b>76</b>	<0.26	<b>350</b>	<28	<b>100</b>	<b>300</b>	<b>53</b>	<110			
<b>Target PAHS</b>																			
Acenaphthene	mg/kg	4	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Acenaphthylene	mg/kg	1	600	10	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Anthracene	mg/kg	1000	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.46</b>	<0.28	<0.78	<b>0.92</b>	<0.22	<1.1			
Benzo(A)Anthracene	mg/kg	7	7	7	<b>0.42</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.79</b>	<0.28	<b>0.78</b>	<b>4</b>	<b>0.59</b>	<1.1			
Benzo(A)Pyrene	mg/kg	2	2	2	<b>0.40</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.70</b>	<0.28	<b>0.91</b>	<b>4.4</b>	<b>0.63</b>	<1.1			
Benzo(B)Fluoranthene	mg/kg	7	7	7	<b>0.35</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.64</b>	<0.28	<b>0.89</b>	<b>4.4</b>	<b>0.57</b>	<1.1			
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<b>0.46</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.51</b>	<0.28	<0.78	<b>3.2</b>	<b>0.42</b>	<1.1			
Benzo(K)Fluoranthene	mg/kg	70	70	70	<b>0.39</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.57</b>	<0.28	<0.78	<b>3.7</b>	<b>0.53</b>	<1.1			
Chrysene	mg/kg	70	70	70	<b>0.52</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.87</b>	<0.28	<b>0.96</b>	<b>5.3</b>	<b>0.74</b>	<1.1			
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Fluoranthene	mg/kg	1000	1000	1000	<b>0.81</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>1.9</b>	<0.28	<b>1.5</b>	<b>9.2</b>	<b>1.3</b>	<1.1			
Fluorene	mg/kg	1000	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.37</b>	<0.28	<0.78	<0.83	<0.22	<1.1			
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<b>0.36</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.36</b>	<0.28	<0.78	<b>2.9</b>	<b>0.36</b>	<1.1			
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.57</b>	<0.28	<0.78	<0.83	<0.22	<1.1			
Naphthalene	mg/kg	4	20	500	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.50</b>	<0.28	<0.78	<b>0.86</b>	<0.22	<1.1			
Phenanthrene	mg/kg	10	500	500	<b>0.44</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>2.1</b>	<0.28	<b>0.80</b>	<b>5.7</b>	<b>1.1</b>	<1.1			
Pyrene	mg/kg	1000	1000	1000	<b>0.88</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>2.1</b>	<0.28	<b>1.7</b>	<b>9</b>	<b>1.4</b>	<1.1			
<b>VPH</b>																			
C5-C8 Aliphatics	mg/kg	100	100	100	NT	NT	<5.5	<b>7.6</b>	<b>26</b>	NT	<b>30</b>	NT	<29	<29	NT	<42			
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	NT	<5.5	<5.4	<b>22</b>	NT	<b>24</b>	NT	<29	<29	NT	<42			
C9-C10 Aromatics	mg/kg	100	100	100	NT	NT	<5.5	<5.4	<b>39</b>	NT	<b>67</b>	NT	<29	<29	NT	<42			
<b>Target VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.30</b>	NT	<0.59	<0.59	NT	<0.83			
Ethylbenzene	mg/kg	40	500	500	NT	NT	<0.11	<0.11	<b>0.48</b>	NT	<b>0.49</b>	NT	<0.59	<0.59	NT	<0.83			
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	NT	<0.11	<0.11	<0.094	NT	<0.11	NT	<0.59	<0.59	NT	<0.83			
Naphthalene	mg/kg	4	20	500	NT	NT	<0.28	<0.27	<b>1.3</b>	NT	<b>1.9</b>	NT	<1.5	<1.5	NT	<2.1			
Toluene	mg/kg	30	500	500	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.42</b>	NT	<0.59	<0.59	NT	<0.83			
m-p Xylene	mg/kg	100	100	500	NT	NT	<0.11	<0.11	<b>1.6</b>	NT	<b>2.7</b>	NT	<0.59	<0.59	NT	<0.83			
o-Xylene	mg/kg	100	100	500	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.45</b>	NT	<0.59	<0.59	NT	<0.83			
<b>Metals</b>																			
Antimony	mg/kg	20	20	20	<b>0.9</b>	NT	<b>0.78</b>	<b>55</b>	<b>5.9</b>	<b>51</b>	<b>10</b>	<b>64</b>	<b>3.5</b>	<b>2.1</b>	NT	<b>3.5</b>			
Arsenic	mg/kg	20	20	20	<b>4</b>	NT	<b>3.2</b>	<b>8.0</b>	<b>3.5</b>	<b>8.1</b>	<b>7.1</b>	<b>23</b>	<b>15</b>	<b>20</b>	<0.51	<b>2.9</b>			
Barium	mg/kg	1000	1000	1000	<b>200</b>	NT	<b>32</b>	<b>280</b>	<b>48</b>	<b>430</b>	<b>470</b>	<b>300</b>	<b>78</b>	<b>86</b>	<b>41</b>	<b>160</b>			
Beryllium	mg/kg	90	90	90	<0.56	NT	<0.59	<0.63	<0.49	<0.65	<0.60	<0.67	<2.0	<2.0	<0.51	<2.8			
Cadmium	mg/kg	70	70	70	<0.56	NT	<0.59	<b>4.4</b>	<b>2.0</b>	<b>5.8</b>	<b>11</b>	<b>3.6</b>	<2.0	<2.0	<0.51	<2.8			
Chromium (III)	mg/kg	1000	1000	1000	NT	NT	NT	<b>33</b>	NT	NT	<b>25</b>	NT	<20	NT	NT	<28			
Chromium (VI)	mg/kg	100	100	100	<b>16</b>	<b>0.50</b>	<b>23</b>	<0.51	<b>19</b>	<b>70</b>	<0.48	<b>64</b>	<1.6	<20	<b>10</b>	<2.2			
Lead	mg/kg	200	200	200	<b>120</b>	NT	<b>150</b>	<b>1200</b>	<b>300</b>	<b>1200</b>	<b>970</b>	<b>1300</b>	<b>180</b>	<b>250</b>	<b>8.1</b>	<b>280</b>			
Mercury	mg/kg	20	20	20	<b>0.17</b>	NT	<0.15	<b>0.60</b>	<0.14	<b>0.18</b>	<b>0.47</b>	<0.19	<0.58	<b>0.99</b>	<0.16	<0.78			
Nickel	mg/kg	600	600	600	<b>12</b>	NT	<b>9.8</b>	<b>76</b>	<b>16</b>	<b>88</b>	<b>40</b>	<b>100</b>	<20	<20	<b>12</b>	<28			
Selenium	mg/kg	400	400	400	<5.36	NT	<5.9	<6.3	<4.9	<6.5	<6.0	<6.7	<20	<20	<5.1	<28			
Silver	mg/kg	100	100	100	<2.8	NT	<3.0	<b>32</b>	<2.5	<b>7.3</b>	<b>12</b>	<9.9	<10	<10	<2.5	<14			
Thallium	mg/kg	8	8	8	<0.56	NT	<0.59	<0.63	<0.49	<0.65	<0.60	<0.67	<2.0	<2.0	<0.51	<2.8			
Vanadium	mg/kg	400	400	400	<b>17</b>	NT	<b>12</b>	<b>20</b>	<b>12</b>	<b>10</b>	<b>18</b>	<b>8.1</b>	<20	<20	<b>16</b>	<28			
Zinc	mg/kg	1000	1000	1000	<b>190</b>	NT	<b>91</b>	<b>2900</b>	<b>250</b>	<b>2900</b>	<b>2400</b>	<b>4000</b>	<b>270</b>	<b>250</b>	<b>19</b>	<b>230</b>			
<b>VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	<b>0.19</b>	NT	NT	NT	NT	NT			
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	<b>0.26</b>	NT	NT	NT	NT	NT			
Cis-1,2-Dichloroethylene	mg/kg	0.1	0.1	100	NT	NT	NT	NT	NT	NT	<b>0.20</b>	NT	NT	NT	NT	NT			
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	<b>0.59</b>	NT	NT	NT	NT	NT			
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	<b>0.34</b>	NT	NT	NT	NT	NT			
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	<b>0.30</b>	NT	NT	NT	NT	NT			
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	<b>1.8</b>	NT	NT	NT	NT	NT			
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	<b>0.93</b>	NT	NT	NT	NT	NT			
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	<b>0.46</b>	NT	NT	NT	NT	NT			
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	<b>0.36</b>	NT	NT	NT	NT	NT			
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	<b>6.9</b>	NT	NT	NT	NT	NT			
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	<b>2.3</b>	NT	NT	NT	NT	NT			
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	<b>3.43</b>	NT	NT	NT	NT	NT			

QC by JMR 5/8/2020

**Abbreviations:**  
EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
PAH = Polycyclic Aromatic Hydrocarbons  
NT = Not Tested  
BDL = Below Detection Limit  
MCP = Massachusetts Contingency Plan  
mg/kg = milligram per kilogram

**Notes:**  
~ = No standard available  
< = indicates parameter not detected above laboratory method reporting limit, shown  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter equals or exceeds the applicable MCP RCS-1  
**BOLD** Parameter exceeds the applicable MCP Method 1 S-1 Cleanup Standard  
1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.

**Table 1**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**618R Waverly Street**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)		Sample Location, Depth, and Date														
					RCS-1	S-1/GW-2	S-1/GW-3	SB-111			SB-112		SB-113			SB-114		SB-115	
								0-1 feet	0-3 feet	4-7 feet	0-3 Feet	5-8 Feet	0-3 feet	5-8 feet	DUP-2	0-3 feet	5-8 feet	0-3 feet	5-8 feet
								4/8/2020	4/8/2020	4/8/2020	4/7/2020	4/7/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020
<b>EPH</b>																			
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<20	<19	<97	<21	<38	<97	<19	<20	<19	<19	<19	<19	<20		
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>85</b>	<b>40</b>	<b>340</b>	<b>50</b>	<b>190</b>	<b>450</b>	<19	<20	<b>40</b>	<b>55</b>	<19	<19	<b>57</b>		
C11-C22 Aromatics	mg/kg	1000	1000	1000	<b>62</b>	<19	<b>200</b>	<b>22</b>	<b>140</b>	<b>97</b>	<19	<20	<b>220</b>	<b>33</b>	<19	<19	<b>75</b>		
<b>Target PAHs</b>																			
Acenaphthene	mg/kg	4	1000	1000	<0.20	<0.19	<b>0.24</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<0.20		
Acenaphthylene	mg/kg	1	600	10	<0.20	<0.19	<0.19	<0.21	<0.38	<0.19	<0.19	<0.20	<b>0.77</b>	<0.19	<0.19	<0.19	<0.20		
Anthracene	mg/kg	1000	1000	1000	<0.20	<0.19	<b>1.3</b>	<0.21	<b>0.84</b>	<0.19	<0.19	<0.20	<b>0.61</b>	<0.19	<0.19	<0.19	<b>0.33</b>		
Benzo(A)Anthracene	mg/kg	7	7	7	<b>1.5</b>	<0.19	<b>4.1</b>	<b>0.39</b>	<b>3</b>	<0.19	<0.19	<0.20	<b>4.9</b>	<b>0.33</b>	<0.19	<0.19	<b>1.6</b>		
Benzo(A)Pyrene	mg/kg	2	2	2	<b>2</b>	<0.19	<b>4.1</b>	<b>0.4</b>	<b>3</b>	<0.19	<0.19	<0.20	<b>4.4</b>	<b>0.37</b>	<0.19	<0.19	<b>1.7</b>		
Benzo(B)Fluoranthene	mg/kg	7	7	7	<b>2.3</b>	<0.19	<b>3.9</b>	<b>0.37</b>	<b>2.8</b>	<0.19	<0.19	<0.20	<b>3.6</b>	<b>0.29</b>	<0.19	<0.19	<b>1.8</b>		
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<b>2</b>	<0.19	<b>3</b>	<b>0.25</b>	<b>2</b>	<0.19	<0.19	<0.20	<b>3.4</b>	<b>0.34</b>	<0.19	<0.19	<b>1.5</b>		
Benzo(K)Fluoranthene	mg/kg	70	70	70	<b>1.7</b>	<0.19	<b>3.4</b>	<b>0.38</b>	<b>2.7</b>	<0.19	<0.19	<0.20	<b>3.4</b>	<b>0.31</b>	<0.19	<0.19	<b>1.4</b>		
Chrysene	mg/kg	70	70	70	<b>1.8</b>	<0.19	<b>4.4</b>	<b>0.42</b>	<b>3.6</b>	<0.19	<0.19	<0.20	<b>5.7</b>	<b>0.38</b>	<0.19	<0.19	<b>1.9</b>		
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<b>0.51</b>	<0.19	<b>0.95</b>	<0.21	<b>0.62</b>	<0.19	<0.19	<0.20	<b>0.78</b>	<0.19	<0.19	<0.19	<b>0.38</b>		
Fluoranthene	mg/kg	1000	1000	1000	<b>2.2</b>	<0.19	<b>8.5</b>	<b>0.8</b>	<b>6.8</b>	<0.19	<0.19	<0.20	<b>9.3</b>	<b>0.77</b>	<0.19	<0.19	<b>3.3</b>		
Fluorene	mg/kg	1000	1000	1000	<0.20	<0.19	<b>0.43</b>	<0.21	<b>0.43</b>	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<b>0.2</b>		
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<b>1.7</b>	<0.19	<b>2.7</b>	<b>0.21</b>	<b>1.7</b>	<0.19	<0.19	<0.20	<b>2.7</b>	<b>0.25</b>	<0.19	<0.19	<b>1.2</b>		
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.20	<0.19	<b>0.22</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<0.20		
Naphthalene	mg/kg	4	20	500	<0.20	<0.19	<b>0.35</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<b>0.26</b>		
Phenanthrene	mg/kg	10	500	500	<b>0.57</b>	<0.19	<b>4.7</b>	<b>0.41</b>	<b>4.7</b>	<0.19	<0.19	<0.20	<b>4.1</b>	<b>0.54</b>	<0.19	<0.19	<b>2.2</b>		
Pyrene	mg/kg	1000	1000	1000	<b>1.9</b>	<0.19	<b>6.7</b>	<b>0.73</b>	<b>5.8</b>	<0.19	<0.19	<0.20	<b>11</b>	<b>0.69</b>	<0.19	<0.19	<b>3.5</b>		
<b>VPH</b>																			
C5-C8 Aliphatics	mg/kg	100	100	100	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3		
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3		
C9-C10 Aromatics	mg/kg	100	100	100	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3		
<b>Target VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
Ethylbenzene	mg/kg	40	500	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
Naphthalene	mg/kg	4	20	500	NT	NT	<b>0.3</b>	NT	<0.52	NT	<0.23	<0.23	<0.20	NT	NT	NT	<0.37		
Toluene	mg/kg	30	500	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
m+p Xylene	mg/kg	100	100	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
o-Xylene	mg/kg	100	100	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15		
<b>Metals</b>																			
Antimony	mg/kg	20	20	20	<b>0.45</b>	<b>29</b>	<b>4.6</b>	<0.50	<b>4.2</b>	<0.56	<0.42	<0.44	<0.48	<b>0.66</b>	<0.40	<0.40	<b>1.7</b>		
Arsenic	mg/kg	20	20	20	<1.7	<b>12</b>	<b>8.2</b>	<b>5.0</b>	<b>16</b>	<b>3.7</b>	<b>2.1</b>	<2.2	<2.4	<15	<b>2.6</b>	<b>8.3</b>			
Barium	mg/kg	1000	1000	1000	<b>51</b>	<b>260</b>	<b>80</b>	<b>38</b>	<b>130</b>	<b>48</b>	<b>26</b>	<b>26</b>	<b>17</b>	<b>210</b>	<b>30</b>	<b>94</b>			
Beryllium	mg/kg	90	90	90	<0.34	<0.48	<b>0.44</b>	<0.50	<0.95	<b>0.69</b>	<0.42	<0.044	<0.48	<0.59	<b>0.56</b>	<0.64			
Cadmium	mg/kg	70	70	70	<0.34	<b>3.8</b>	<b>0.7</b>	<0.50	<b>1.1</b>	<0.56	<0.42	<0.44	<0.48	<0.59	<0.40	<0.64			
Chromium (III)	mg/kg	1000	1000	1000	NT	NT	<b>13</b>	NT	<b>21</b>	NT	<b>5.7</b>	NT	NT	<29	NT	<b>15</b>			
Chromium (VI)	mg/kg	100	100	100	<b>12</b>	<b>43</b>	<b>2</b>	<b>18</b>	<b>1.3</b>	<b>15</b>	<0.40	<b>6.3</b>	<4.8	<0.59	<b>8</b>	<0.59			
Lead	mg/kg	200	200	200	<b>21</b>	<b>630</b>	<b>460</b>	<b>66</b>	<b>470</b>	<b>36</b>	<b>3.2</b>	<b>3.4</b>	<b>21</b>	<b>270</b>	<b>14</b>	<b>330</b>			
Mercury	mg/kg	20	20	20	<0.14	<b>0.26</b>	<0.16	<0.15	<b>0.80</b>	<0.16	<0.15	<0.14	<0.15	<0.21	<0.13	<0.21			
Nickel	mg/kg	600	600	600	<b>11</b>	<b>52</b>	<b>18</b>	<b>18</b>	<b>17</b>	<5.6	<b>6.2</b>	<b>5.9</b>	<4.8	<29	<b>5.4</b>	<b>9.9</b>			
Selenium	mg/kg	400	400	400	<3.4	<4.8	<4.0	<5.0	<9.5	<5.6	<4.2	<4.4	<4.8	<29	<4.0	<6.4			
Silver	mg/kg	100	100	100	<1.7	<b>4.5</b>	<2.0	<2.5	<4.8	<2.8	<2.1	<2.4	<2.9	<2.0	<2.0	<3.2			
Thallium	mg/kg	8	8	8	<0.34	<b>0.48</b>	<b>0.4</b>	<0.50	<0.95	<0.56	<0.42	<0.44	<0.48	<0.59	<0.40	<0.64			
Vanadium	mg/kg	400	400	400	<b>19</b>	<b>12</b>	<b>14</b>	<b>21</b>	<b>26</b>	<b>16</b>	<b>7.8</b>	<b>8.3</b>	<4.8	<29	<b>11</b>	<b>16</b>			
Zinc	mg/kg	1000	1000	1000	<b>42</b>	<b>1500</b>	<b>330</b>	<b>51</b>	<b>300</b>	<b>43</b>	<b>18</b>	<b>18</b>	<b>40</b>	<b>340</b>	<b>28</b>	<b>210</b>			
<b>VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Cis-1,2-Dichloroethylene	mg/kg	0.1	100	100	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT		

QC by JMR 5/8/2020

**Abbreviations:**  
EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
PAH = Polycyclic Aromatic Hydrocarbons  
NT = Not Tested  
BDL = Below Detection Limit  
MCP = Massachusetts Contingency Plan  
mg/kg = milligram per kilogram

**Notes:**  
~ = No standard available  
< = indicates parameter not detected above laboratory method reporting limit, shown  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter equals or exceeds the applicable MCP RCS-1  
**BOLD** Parameter exceeds the applicable MCP Method 1 S-1 Cleanup Standard  
1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.

Table 2  
 Summary of Soil Analytical Results - Shallow Lead Concentrations  
 618R Waverly Street  
 Framingham, Massachusetts

Parameters	Units	MCP - Method 1 Cleanup Standards <sup>1</sup>	Sample Location, Date, and Depth							
			SB-102	SB-106	SB-107	SB-108	SB-109	SB-110	SB-111	SB-112
			4/6/2020 0-1 feet	4/6/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/8/2020 0-1 feet	4/7/2020 0-1 feet
Metals LEAD	mg/Kg	200	73	68	67	900	790	11	21	25

QC by JRS 7/24/2020

**Abbreviations:**

MCP = Massachusetts Contingency Plan  
 mg/kg = milligram per kilogram

**Notes:**

< = indicates parameter not detected above laboratory method reporting limit, shown  
 1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter exceeds the MCP Method 1, S-1/GW-3 Cleanup Standard

## APPENDIX A

### Public Notification Letters

March 11, 2021

Mr. Sam Wong  
Director, Framingham Health Dept.  
Framingham City Hall  
150 Concord Street  
Framingham, Massachusetts 01702

Re: **Immediate Response Action (IRA) Completion Report**  
**618R Waverly Street, Framingham, MA (Parcel IDs 134-64-7867 and 134-64-9905)**  
**Release Tracking Number (RTN) 3-36304**

Dear Mr. Wong:

Weston & Sampson is hereby notifying your office that an Immediate Response Action (IRA) Completion Report has been submitted to the Massachusetts Department of Environmental Protection (MassDEP) for the above-referenced Release Tracking Number (RTN). The RTN is located on two parcels of land, owned by the City, identified on the Framingham Assessor's maps as Parcel IDs 134-64-7867 and 134-64-9905 (Site). Implementation of an IRA, as outlined in 310 CMR 40.0410, was necessary at the Site to mitigate an Imminent Hazard (IH) due to elevated lead concentrations in shallow soil.

The IRA activities completed at the Site included the installation of temporary chain-link fence, which mitigated the IH condition, and will remain in-place until Comprehensive Response Actions are undertaken. Future response actions will likely include the collection and analysis of shallow soil samples and the excavation of shallow lead-impacted soil.

This notification is provided in accordance with the public involvement requirements of the MCP at 310 CMR 40.1403(3)(c). The IRA Completion Report is available for public review online at the MassDEP's website (<https://eeaonline.eea.state.ma.us/portal#!/wastesite/3-0036304>).

If you have any questions or concerns regarding this notice, please do not hesitate to contact the undersigned at (978) 532-1900.

Sincerely,

WESTON & SAMPSON ENGINEERS, INC.



Frank Ricciardi, PE, LSP  
Vice President

Cc: Mayor Yvonne M. Spicer, City of Framingham  
MassDEP, Bureau of Waste Site Cleanup

March 11, 2021

Honorable Mayor Yvonne M. Spicer  
Framingham City Hall  
150 Concord Street  
Framingham, Massachusetts 01702

Re: **Immediate Response Action (IRA) Completion Report**  
**618R Waverly Street, Framingham, MA (Parcel IDs 134-64-7867 and 134-64-9905)**  
**Release Tracking Number (RTN) 3-36304**

Dear Mayor Spicer:

Weston & Sampson is hereby notifying your office that an Immediate Response Action (IRA) Completion Report has been submitted to the Massachusetts Department of Environmental Protection (MassDEP) for the above-referenced Release Tracking Number (RTN). The RTN is located on two parcels of land, owned by the City, identified on the Framingham Assessor's maps as Parcel IDs 134-64-7867 and 134-64-9905 (Site). Implementation of an IRA, as outlined in 310 CMR 40.0410, was necessary at the Site to mitigate an Imminent Hazard (IH) due to elevated lead concentrations in shallow soil.

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This notification is provided in accordance with the public involvement requirements of the MCP at 310 CMR 40.1403(3)(c). The IRA Completion Report is available for public review online at the MassDEP's website (<https://eeaonline.eea.state.ma.us/portal#!/wastesite/3-0036304>).

If you have any questions or concerns regarding this notice, please do not hesitate to contact the undersigned at (978) 532-1900.

Sincerely,

WESTON & SAMPSON ENGINEERS, INC.



Frank Ricciardi, PE, LSP  
Vice President

Cc: Mr. Sam Wong, Director, Framingham Health Dept.  
MassDEP, Bureau of Waste Site Cleanup



## APPENDIX B

### Laboratory Analytical Reports

# Laboratory Report



**Absolute Resource** *associates*

124 Heritage Avenue Portsmouth NH 03801

Jill Murphy  
Weston & Sampson  
55 Walkers Brook Drive  
Reading, MA 01867

PO Number: Framingham Brownfields  
Job ID: 52498  
Date Received: 4/6/20

Project: Cedar Woods 2180311

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below. The reported results apply to the sample(s) in the condition as received at the time the laboratory took custody. This report shall not be reproduced except in full and with approval from the laboratory. The liability of ARA is limited to the cost of the requested analyses, unless otherwise agreed upon in writing.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,  
Absolute Resource Associates

A handwritten signature in black ink, appearing to read 'A. DeWees', written in a cursive style.

Aaron DeWees  
Chief Operating Officer

Date of Approval: 5/4/2020  
Total number of pages: 48

## Absolute Resource Associates Certifications

New Hampshire 1732  
Maine NH902

Massachusetts M-NH902

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-106 (0-3')	Solid	4/6/2020 9:40	52498-002	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-106 (5-8')	Solid	4/6/2020 10:00	52498-003	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
SB-106 (3-4')	Solid	4/6/2020 9:50	52498-004	Hexavalent Chromium in solids by SW3060A7196A ORP in solids by ASTM-1498-08 pH in solids by SW9045C

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-105 (0-3')	Solid	4/6/2020 10:40	52498-005	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-105 (5-8')	Solid	4/6/2020 11:00	52498-006	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-102 (0-3')	Solid	4/6/2020 11:45	52498-007	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-102 (11-14')	Solid	4/6/2020 11:40	52498-009	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-101 (0-3')	Solid	4/6/2020 12:55	52498-010	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-101 (9-12')	Solid	4/6/2020 13:05	52498-011	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-103 (0-3')	Solid	4/6/2020 12:20	52498-012	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-103 (3-5')	Solid	4/6/2020 12:20	52498-013	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-103 (7-10')	Solid	4/6/2020 12:25	52498-014	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
Trip Blank	Solid	4/6/2020 0:00	52498-015	VPH in solids by MA DEP Method
SB-106 (5-8')REX	Solid	4/6/2020 10:00	52498-016	EPH in solids by MADEP Method



Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-003

Sample ID: SB-106 (5-8')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.7 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 10:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 5.5	5.5	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.5	5.5	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
benzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
toluene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
ethylbenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
m&p-xylenes	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
o-xylene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
naphthalene	< 0.28	0.28	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
C5-C8 Aliphatics	< 5.5	5.5	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
C9-C12 Aliphatics	< 5.5	5.5	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
C9-C10 Aromatics	< 5.5	5.5	ug/g	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>108</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>108</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>91</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	12:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-006

Sample ID: SB-105 (5-8')

Matrix: Solid Percent Dry: 88% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.65 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 11:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 4.3	4.3	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 4.3	4.3	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
benzene	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
toluene	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
ethylbenzene	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
m&p-xylenes	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
o-xylene	< 0.087	0.087	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
naphthalene	< 0.22	0.22	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
C5-C8 Aliphatics	< 4.3	4.3	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
C9-C12 Aliphatics	< 4.3	4.3	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
C9-C10 Aromatics	< 4.3	4.3	ug/g	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>107</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>107</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>99</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-009

Sample ID: SB-102 (11-14')

Matrix: Solid Percent Dry: 40% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.65 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 11:40

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 16	16	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 16	16	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
benzene	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
toluene	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
ethylbenzene	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
m&p-xylenes	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
o-xylene	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
naphthalene	< 0.78	0.78	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
C5-C8 Aliphatics	< 16	16	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
C9-C12 Aliphatics	< 16	16	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
C9-C10 Aromatics	< 16	16	ug/g	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>110</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>110</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>106</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	13:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-011

Sample ID: SB-101 (9-12')

Matrix: Solid Percent Dry: 74% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 13:05

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 6.3	6.3	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 6.3	6.3	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
benzene	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
toluene	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
ethylbenzene	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
m&p-xylenes	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
o-xylene	< 0.13	0.13	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
naphthalene	< 0.31	0.31	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
C5-C8 Aliphatics	< 6.3	6.3	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
C9-C12 Aliphatics	< 6.3	6.3	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
C9-C10 Aromatics	< 6.3	6.3	ug/g	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>102</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>103</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>69 *</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:19	MA VPH

\* **This surrogate showed recovery outside the acceptance limits.**

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-014

Sample ID: SB-103 (7-10')

Matrix: Solid Percent Dry: 76.8% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.66 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 12:25

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 5.8	5.8	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.8	5.8	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
benzene	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
toluene	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
ethylbenzene	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
m&p-xylenes	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
o-xylene	< 0.12	0.12	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
naphthalene	< 0.29	0.29	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
C5-C8 Aliphatics	< 5.8	5.8	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
C9-C12 Aliphatics	< 5.8	5.8	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
C9-C10 Aromatics	< 5.8	5.8	ug/g	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>104</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>105</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>78</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	14:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-015

Sample ID: Trip Blank

Matrix: Solid

Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil

Received on ice at 0°C, in satisfactory condition.

Sampled: 4/6/20 0:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
benzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
toluene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
ethylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
m&p-xylenes	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
o-xylene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
naphthalene	< 0.25	0.25	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
C5-C8 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
C9-C12 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
C9-C10 Aromatics	< 5.0	5.0	ug/g	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>105</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>105</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>90</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	11:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-002

Sample ID: SB-106 (0-3')

Matrix: Solid Percent Dry: 89.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 9:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
2-methylnaphthalene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
phenanthrene	<b>0.44</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
acenaphthene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
acenaphthylene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
fluorene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
anthracene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
fluoranthene	<b>0.81</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
pyrene	<b>0.88</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
benzo(a)anthracene	<b>0.42</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
chrysene	<b>0.52</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
benzo(b)fluoranthene	<b>0.35</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
benzo(k)fluoranthene	<b>0.39</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
benzo(a)pyrene	<b>0.40</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
indeno(1,2,3-cd)pyrene	<b>0.36</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
dibenzo(a,h)anthracene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
benzo(g,h,i)perylene	<b>0.46</b>	0.21	ug/g	1	CL	4/7/20	12625	4/9/20	7:07	MA EPH
Unadjusted C11-C22 Aromatics	<b>49</b>	21	ug/g	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
C9-C18 Aliphatics	< 21	21	ug/g	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
C19-C36 Aliphatics	<b>91</b>	21	ug/g	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
C11-C22 Aromatics	<b>44</b>	21	ug/g	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>55</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
o-terphenyl SUR	<b>56</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
2-fluorobiphenyl SUR	<b>66</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH
2-bromonaphthalene SUR	<b>62</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	22:29	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-003

Sample ID: SB-106 (5-8')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
2-methylnaphthalene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
phenanthrene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
acenaphthene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
acenaphthylene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
fluorene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
anthracene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
fluoranthene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
pyrene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
benzo(a)anthracene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
chrysene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
benzo(b)fluoranthene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
benzo(k)fluoranthene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
benzo(a)pyrene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
indeno(1,2,3-cd)pyrene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
dibenzo(a,h)anthracene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
benzo(g,h,i)perylene	< 0.47	0.47	ug/g	1	CL	4/7/20	12625	4/9/20	8:07	MA EPH
Unadjusted C11-C22 Aromatics	<b>290</b>	240	ug/g	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
C9-C18 Aliphatics	< 240	240	ug/g	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
C19-C36 Aliphatics	<b>570</b>	240	ug/g	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
C11-C22 Aromatics	<b>290</b>	240	ug/g	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>17 *</b>	40-140	%	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
o-terphenyl SUR	<b>17 *</b>	40-140	%	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
2-fluorobiphenyl SUR	<b>141 *</b>	40-140	%	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH
2-bromonaphthalene SUR	<b>125</b>	40-140	%	5	DBV	4/7/20	12625	4/8/20	17:17	MA EPH

\* The surrogate showed recovery outside the acceptance limits. Re-extraction beyond method holding time produced similar results. The re-extracted results are included.



Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-005

Sample ID: SB-105 (0-3')

Matrix: Solid Percent Dry: 81.3% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
2-methylnaphthalene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
phenanthrene	1.5	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
acenaphthene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
acenaphthylene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
fluorene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
anthracene	< 0.24	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
fluoranthene	3.2	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
pyrene	2.9	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
benzo(a)anthracene	1.5	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
chrysene	1.9	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
benzo(b)fluoranthene	1.6	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
benzo(k)fluoranthene	1.2	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
benzo(a)pyrene	1.4	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
indeno(1,2,3-cd)pyrene	1.1	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
dibenzo(a,h)anthracene	0.38	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
benzo(g,h,i)perylene	1.2	0.24	ug/g	1	CL	4/7/20	12625	4/9/20	3:36	MA EPH
Unadjusted C11-C22 Aromatics	150	24	ug/g	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
C9-C18 Aliphatics	< 24	24	ug/g	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
C19-C36 Aliphatics	280	24	ug/g	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
C11-C22 Aromatics	130	24	ug/g	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	47	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
o-terphenyl SUR	49	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
2-fluorobiphenyl SUR	69	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH
2-bromonaphthalene SUR	67	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:34	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-006

Sample ID: SB-105 (5-8')

Matrix: Solid Percent Dry: 88% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
2-methylnaphthalene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
phenanthrene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
acenaphthene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
acenaphthylene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
fluorene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
anthracene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
fluoranthene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
pyrene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
benzo(a)anthracene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
chrysene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
benzo(b)fluoranthene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
benzo(k)fluoranthene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
benzo(a)pyrene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
indeno(1,2,3-cd)pyrene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
dibenzo(a,h)anthracene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
benzo(g,h,i)perylene	< 0.22	0.22	ug/g	1	CL	4/7/20	12625	4/8/20	18:33	MA EPH
Unadjusted C11-C22 Aromatics	< 22	22	ug/g	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
C9-C18 Aliphatics	< 22	22	ug/g	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
C19-C36 Aliphatics	<b>23</b>	22	ug/g	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
C11-C22 Aromatics	< 22	22	ug/g	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>58</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
o-terphenyl SUR	<b>58</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
2-fluorobiphenyl SUR	<b>66</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH
2-bromonaphthalene SUR	<b>64</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:18	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-007

Sample ID: SB-102 (0-3')

Matrix: Solid Percent Dry: 93% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
2-methylnaphthalene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
phenanthrene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
acenaphthene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
acenaphthylene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
fluorene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
anthracene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
fluoranthene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
pyrene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
benzo(a)anthracene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
chrysene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
benzo(b)fluoranthene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
benzo(k)fluoranthene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
benzo(a)pyrene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
indeno(1,2,3-cd)pyrene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
dibenzo(a,h)anthracene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
benzo(g,h,i)perylene	< 0.20	0.20	ug/g	1	CL	4/7/20	12625	4/9/20	7:37	MA EPH
Unadjusted C11-C22 Aromatics	<b>51</b>	20	ug/g	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
C9-C18 Aliphatics	< 20	20	ug/g	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
C19-C36 Aliphatics	<b>130</b>	20	ug/g	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
C11-C22 Aromatics	<b>50</b>	20	ug/g	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>43</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
o-terphenyl SUR	<b>44</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
2-fluorobiphenyl SUR	<b>66</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH
2-bromonaphthalene SUR	<b>62</b>	40-140	%	1	DBV	4/7/20	12625	4/9/20	23:03	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-009

Sample ID: SB-102 (11-14')

Matrix: Solid Percent Dry: 40% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
2-methylnaphthalene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
phenanthrene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
acenaphthene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
acenaphthylene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
fluorene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
anthracene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
fluoranthene	<b>0.48</b>	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
pyrene	<b>0.56</b>	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
benzo(a)anthracene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
chrysene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
benzo(b)fluoranthene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
benzo(k)fluoranthene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
benzo(a)pyrene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
indeno(1,2,3-cd)pyrene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
dibenzo(a,h)anthracene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
benzo(g,h,i)perylene	< 0.46	0.46	ug/g	1	CL	4/7/20	12625	4/9/20	2:36	MA EPH
Unadjusted C11-C22 Aromatics	< 46	46	ug/g	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
C9-C18 Aliphatics	< 46	46	ug/g	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
C19-C36 Aliphatics	<b>80</b>	46	ug/g	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
C11-C22 Aromatics	< 46	46	ug/g	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>54</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
o-terphenyl SUR	<b>58</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
2-fluorobiphenyl SUR	<b>66</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH
2-bromonaphthalene SUR	<b>62</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	15:00	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-010

Sample ID: SB-101 (0-3')

Matrix: Solid Percent Dry: 90.7% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
2-methylnaphthalene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
phenanthrene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
acenaphthene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
acenaphthylene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
fluorene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
anthracene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
fluoranthene	<b>0.22</b>	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
pyrene	<b>0.24</b>	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
benzo(a)anthracene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
chrysene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
benzo(b)fluoranthene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
benzo(k)fluoranthene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
benzo(a)pyrene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
indeno(1,2,3-cd)pyrene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
dibenzo(a,h)anthracene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
benzo(g,h,i)perylene	< 0.21	0.21	ug/g	1	CL	4/7/20	12625	4/8/20	22:35	MA EPH
Unadjusted C11-C22 Aromatics	< 21	21	ug/g	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
C9-C18 Aliphatics	< 21	21	ug/g	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
C19-C36 Aliphatics	<b>35</b>	21	ug/g	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
C11-C22 Aromatics	< 21	21	ug/g	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>61</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
o-terphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
2-fluorobiphenyl SUR	<b>67</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH
2-bromonaphthalene SUR	<b>64</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	13:52	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-011

Sample ID: SB-101 (9-12')

Matrix: Solid Percent Dry: 74% Results expressed on a dry weight basis.

Sampled: 4/6/20 13:05

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
2-methylnaphthalene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
phenanthrene	<b>2.1</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
acenaphthene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
acenaphthylene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
fluorene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
anthracene	<b>0.40</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
fluoranthene	<b>2.1</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
pyrene	<b>2.3</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
benzo(a)anthracene	<b>0.98</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
chrysene	<b>1.0</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
benzo(b)fluoranthene	<b>0.88</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
benzo(k)fluoranthene	<b>0.71</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
benzo(a)pyrene	<b>0.89</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
indeno(1,2,3-cd)pyrene	<b>0.63</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
dibenzo(a,h)anthracene	< 0.25	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
benzo(g,h,i)perylene	<b>0.72</b>	0.25	ug/g	1	CL	4/7/20	12625	4/9/20	3:06	MA EPH
Unadjusted C11-C22 Aromatics	<b>54</b>	25	ug/g	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
C9-C18 Aliphatics	< 25	25	ug/g	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
C19-C36 Aliphatics	<b>82</b>	25	ug/g	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
C11-C22 Aromatics	<b>41</b>	25	ug/g	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>57</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
o-terphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
2-fluorobiphenyl SUR	<b>69</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH
2-bromonaphthalene SUR	<b>67</b>	40-140	%	1	DBV	4/7/20	12625	4/8/20	14:26	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-016

Sample ID: SB-106 (5-8')REX

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
2-methylnaphthalene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
phenanthrene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
acenaphthene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
acenaphthylene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
fluorene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
anthracene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
fluoranthene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
pyrene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
benzo(a)anthracene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
chrysene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
benzo(b)fluoranthene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
benzo(k)fluoranthene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
benzo(a)pyrene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
indeno(1,2,3-cd)pyrene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
dibenzo(a,h)anthracene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
benzo(g,h,i)perylene	< 2.4	2.4	ug/g	5	CL	4/9/20	12630	4/9/20	21:15	MA EPH
Unadjusted C11-C22 Aromatics	<b>410</b>	240	ug/g	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
C9-C18 Aliphatics	< 240	240	ug/g	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
C19-C36 Aliphatics	<b>830</b>	240	ug/g	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
C11-C22 Aromatics	<b>410</b>	240	ug/g	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>21</b> *	40-140	%	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
o-terphenyl SUR	<b>23</b> *	40-140	%	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
2-fluorobiphenyl SUR	<b>118</b>	40-140	%	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH
2-bromonaphthalene SUR	<b>107</b>	40-140	%	5	DBV	4/9/20	12630	4/9/20	23:38	MA EPH

\* The surrogate showed recovery outside the acceptance limits.

Note: Dilution was required due to sample matrix interference.

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-002

Sample ID: SB-106 (0-3')

Matrix: Solid Percent Dry: 89.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 9:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	0.90	0.56	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
Arsenic	4.0	2.8	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Barium	200 DM	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
D = The RPD for the sample duplicate, run as internal QC, was outside the 20% acceptance range. M = The recovery for the matrix spike was 161%. The acceptance criteria is 75-125%.										
Beryllium	< 0.56	0.56	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Cadmium	< 0.56	0.56	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Chromium	16	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
Lead	120	2.8	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
Mercury	0.17	0.14	ug/g	1	AGN	4/7/20	12622	4/7/20	18:25	SW7471B
Nickel	12	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
Selenium	< 5.6	5.6	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Silver	< 2.8	2.8	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Thallium	< 0.56	0.56	ug/g	5	AGN	4/8/20	12626	4/9/20	2:01	SW3051A6020A
Vanadium	17	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A
Zinc	190 M	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	0:52	SW3051A6020A

M = The recovery for the matrix spike was 141%. The acceptance criteria is 75-125%.

Sample#: 52498-003

Sample ID: SB-106 (5-8')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	0.78	0.59	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Arsenic	3.2	3.0	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Barium	32	5.9	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Beryllium	< 0.59	0.59	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Cadmium	< 0.59	0.59	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Chromium	23	5.9	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Lead	150	3.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/7/20	12622	4/7/20	18:27	SW7471B
Nickel	9.8	5.9	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Selenium	< 5.9	5.9	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Silver	< 3.0	3.0	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Thallium	< 0.59	0.59	ug/g	5	AGN	4/8/20	12626	4/9/20	2:29	SW3051A6020A
Vanadium	12	5.9	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A
Zinc	91	5.9	ug/g	5	AGN	4/8/20	12626	4/10/20	1:20	SW3051A6020A



Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-005

Sample ID: SB-105 (0-3')

Matrix: Solid Percent Dry: 81.3% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	4.7	0.60	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Arsenic	7.3	3.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Barium	310	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Beryllium	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	2:38	SW3051A6020A
Cadmium	2.0	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	2:38	SW3051A6020A
Chromium	18	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Lead	650	3.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Mercury	2.9	0.79	ug/g	5	AGN	4/7/20	12622	4/7/20	18:53	SW7471B
Nickel	18	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Selenium	< 6.0	6.0	ug/g	5	AGN	4/8/20	12626	4/9/20	2:38	SW3051A6020A
Silver	< 3.0	3.0	ug/g	5	AGN	4/8/20	12626	4/9/20	2:38	SW3051A6020A
Thallium	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	2:38	SW3051A6020A
Vanadium	19	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A
Zinc	770	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	1:29	SW3051A6020A

Sample#: 52498-006

Sample ID: SB-105 (5-8')

Matrix: Solid Percent Dry: 88% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.55	0.55	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Arsenic	3.0	2.7	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Barium	23	5.5	ug/g	5	AGN	4/8/20	12626	4/10/20	1:39	SW3051A6020A
Beryllium	< 0.55	0.55	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Cadmium	< 0.55	0.55	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Chromium	11	5.5	ug/g	5	AGN	4/8/20	12626	4/10/20	1:39	SW3051A6020A
Lead	5.0	2.7	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/7/20	12622	4/7/20	18:31	SW7471B
Nickel	14	5.5	ug/g	5	AGN	4/8/20	12626	4/10/20	1:39	SW3051A6020A
Selenium	< 5.5	5.5	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Silver	< 2.7	2.7	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Thallium	< 0.55	0.55	ug/g	5	AGN	4/8/20	12626	4/9/20	3:16	SW3051A6020A
Vanadium	14	5.5	ug/g	5	AGN	4/8/20	12626	4/10/20	1:39	SW3051A6020A
Zinc	69	5.5	ug/g	5	AGN	4/8/20	12626	4/10/20	1:39	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-007

Sample ID: SB-102 (0-3')

Matrix: Solid Percent Dry: 93% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.54	0.54	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Arsenic	<b>3.3</b>	2.7	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Barium	<b>27</b>	5.4	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A
Beryllium	< 0.54	0.54	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Cadmium	< 0.54	0.54	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Chromium	<b>17</b>	5.4	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A
Lead	<b>16</b>	2.7	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A
Mercury	< 0.13	0.13	ug/g	1	AGN	4/7/20	12622	4/7/20	18:33	SW7471B
Nickel	<b>23</b>	5.4	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A
Selenium	< 5.4	5.4	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Silver	< 2.7	2.7	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Thallium	< 0.54	0.54	ug/g	5	AGN	4/8/20	12626	4/9/20	3:25	SW3051A6020A
Vanadium	<b>20</b>	5.4	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A
Zinc	<b>31</b>	5.4	ug/g	5	AGN	4/8/20	12626	4/10/20	1:48	SW3051A6020A

Sample#: 52498-009

Sample ID: SB-102 (11-14')

Matrix: Solid Percent Dry: 40% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 1.1	1.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Arsenic	<b>10</b>	5.6	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Barium	<b>56</b>	11	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Beryllium	< 1.1	1.1	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Cadmium	< 1.1	1.1	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Chromium	<b>41</b>	11	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Lead	<b>190</b>	5.6	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Mercury	< 0.31	0.31	ug/g	1	AGN	4/7/20	12622	4/7/20	18:34	SW7471B
Nickel	<b>64</b>	11	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Selenium	< 11	11	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Silver	< 5.6	5.6	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Thallium	< 1.1	1.1	ug/g	5	AGN	4/8/20	12626	4/9/20	3:34	SW3051A6020A
Vanadium	<b>17</b>	11	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A
Zinc	<b>530</b>	11	ug/g	5	AGN	4/8/20	12626	4/10/20	2:25	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-010

Sample ID: SB-101 (0-3')

Matrix: Solid Percent Dry: 90.7% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.52	0.52	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Arsenic	<b>3.3</b>	2.6	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Barium	<b>18</b>	5.2	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A
Beryllium	< 0.52	0.52	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Cadmium	< 0.52	0.52	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Chromium	<b>7.3</b>	5.2	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A
Lead	<b>8.7</b>	2.6	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/7/20	12622	4/7/20	18:36	SW7471B
Nickel	<b>7.4</b>	5.2	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A
Selenium	< 5.2	5.2	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Silver	< 2.6	2.6	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Thallium	< 0.52	0.52	ug/g	5	AGN	4/8/20	12626	4/9/20	3:44	SW3051A6020A
Vanadium	<b>11</b>	5.2	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A
Zinc	<b>17</b>	5.2	ug/g	5	AGN	4/8/20	12626	4/10/20	2:35	SW3051A6020A

Sample#: 52498-011

Sample ID: SB-101 (9-12')

Matrix: Solid Percent Dry: 74% Results expressed on a dry weight basis.

Sampled: 4/6/20 13:05

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	<b>2.7</b>	0.68	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Arsenic	<b>28</b>	3.4	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Barium	<b>140</b>	6.8	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Beryllium	< 0.68	0.68	ug/g	5	AGN	4/8/20	12626	4/9/20	3:53	SW3051A6020A
Cadmium	<b>1.3</b>	0.68	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Chromium	<b>51</b>	6.8	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Lead	<b>230</b>	3.4	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Mercury	<b>0.58</b>	0.18	ug/g	1	AGN	4/7/20	12622	4/7/20	18:42	SW7471B
Nickel	<b>14</b>	6.8	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Selenium	< 6.8	6.8	ug/g	5	AGN	4/8/20	12626	4/9/20	3:53	SW3051A6020A
Silver	< 3.4	3.4	ug/g	5	AGN	4/8/20	12626	4/9/20	3:53	SW3051A6020A
Thallium	< 0.68	0.68	ug/g	5	AGN	4/8/20	12626	4/9/20	3:53	SW3051A6020A
Vanadium	<b>15</b>	6.8	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A
Zinc	<b>220</b>	6.8	ug/g	5	AGN	4/8/20	12626	4/10/20	2:44	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-012

Sample ID: SB-103 (0-3')

Matrix: Solid Percent Dry: 93% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.51	0.51	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Arsenic	<b>3.5</b>	2.5	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Barium	<b>25</b>	5.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A
Beryllium	< 0.51	0.51	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Cadmium	< 0.51	0.51	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Chromium	<b>6.9</b>	5.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A
Lead	<b>22</b>	2.5	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/7/20	12622	4/7/20	18:48	SW7471B
Nickel	<b>6.3</b>	5.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A
Selenium	< 5.1	5.1	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Silver	< 2.5	2.5	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Thallium	< 0.51	0.51	ug/g	5	AGN	4/8/20	12626	4/9/20	4:02	SW3051A6020A
Vanadium	<b>10</b>	5.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A
Zinc	<b>19</b>	5.1	ug/g	5	AGN	4/8/20	12626	4/10/20	2:53	SW3051A6020A

Sample#: 52498-013

Sample ID: SB-103 (3-5')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	<b>3.7</b>	0.62	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Arsenic	<b>13</b>	3.1	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Barium	<b>250</b>	6.2	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Beryllium	< 0.62	0.62	ug/g	5	AGN	4/8/20	12626	4/9/20	4:12	SW3051A6020A
Cadmium	<b>4.9</b>	0.62	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Chromium	<b>9.4</b>	6.2	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Lead	<b>710</b>	3.1	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Mercury	<b>0.58</b>	0.16	ug/g	1	AGN	4/7/20	12622	4/7/20	18:49	SW7471B
Nickel	<b>12</b>	6.2	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Selenium	< 6.2	6.2	ug/g	5	AGN	4/8/20	12626	4/9/20	4:12	SW3051A6020A
Silver	< 3.1	3.1	ug/g	5	AGN	4/8/20	12626	4/9/20	4:12	SW3051A6020A
Thallium	< 0.62	0.62	ug/g	5	AGN	4/8/20	12626	4/9/20	4:12	SW3051A6020A
Vanadium	<b>10.0</b>	6.2	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A
Zinc	<b>920</b>	6.2	ug/g	5	AGN	4/8/20	12626	4/10/20	3:03	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-014

Sample ID: SB-103 (7-10')

Matrix: Solid Percent Dry: 76.8% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:25

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Arsenic	<b>3.3</b>	3.0	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Barium	<b>36</b>	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A
Beryllium	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Cadmium	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Chromium	<b>18</b>	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A
Lead	<b>8.4</b>	3.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A
Mercury	< 0.17	0.17	ug/g	1	AGN	4/7/20	12622	4/7/20	18:51	SW7471B
Nickel	<b>13</b>	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A
Selenium	< 6.0	6.0	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Silver	< 3.0	3.0	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Thallium	< 0.60	0.60	ug/g	5	AGN	4/8/20	12626	4/9/20	4:21	SW3051A6020A
Vanadium	<b>25</b>	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A
Zinc	<b>18</b>	6.0	ug/g	5	AGN	4/8/20	12626	4/10/20	3:12	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52498

Sample#: 52498-004

Sample ID: SB-106 (3-4')

Matrix: Solid Percent Dry: 88.6% Results expressed on a dry weight basis.

Sampled: 4/6/20 9:50

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	0.50	0.45	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	430		mV	1	WAS	2001824	4/7/20	8:00	ASTM1498-08	
pH	5.9		pH	1	WAS	2001823	4/7/20	7:47	SW9045C	

Sample#: 52498-005

Sample ID: SB-105 (0-3')

Matrix: Solid Percent Dry: 81.3% Results expressed on a dry weight basis.

Sampled: 4/6/20 10:40

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.48	0.48	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	450		mV	1	WAS	2001824	4/7/20	8:10	ASTM1498-08	
pH	6.8		pH	1	WAS	2001823	4/7/20	7:53	SW9045C	

Sample#: 52498-009

Sample ID: SB-102 (11-14')

Matrix: Solid Percent Dry: 40% Results expressed on a dry weight basis.

Sampled: 4/6/20 11:40

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 1.00	1.00	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	49		mV	1	WAS	2001824	4/7/20	8:14	ASTM1498-08	
pH	6.9		pH	1	WAS	2001823	4/7/20	8:04	SW9045C	

Sample#: 52498-011

Sample ID: SB-101 (9-12')

Matrix: Solid Percent Dry: 74% Results expressed on a dry weight basis.

Sampled: 4/6/20 13:05

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	3.8	2.8	ug/g	5	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	180		mV	1	WAS	2001824	4/7/20	8:26	ASTM1498-08	
pH	6.1		pH	1	WAS	2001823	4/7/20	7:55	SW9045C	

Sample#: 52498-013

Sample ID: SB-103 (3-5')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/6/20 12:20

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.49	0.49	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	250 D		mV	1	WAS	2001824	4/7/20	8:39	ASTM1498-08	
D = The RPD for the sample duplicate, run as internal QC, was outside the 10mV acceptance range. The duplicate result is 370 mV.										
pH	6.4		pH	1	WAS	2001823	4/7/20	8:00	SW9045C	

# Quality Control Report



124 Heritage Avenue Unit 16  
Portsmouth, NH 03801

[www.absoluteresourceassociates.com](http://www.absoluteresourceassociates.com)

## MassDEP Analytical Protocol Certification Form

Laboratory Name: Absolute Resource Associates

Project #: 2180311

Project Location: Massachusetts

RTN:

**This Form provides certifications for the following data set: list Laboratory Sample ID Number(s): 52498**

Matrices:  Groundwater/Surface Water  Soil/Sediment  Drinking Water  Air  Other:

**CAM Protocol** (check all that apply below):

8260 VOC CAM II A <input type="checkbox"/>	7470/7471 Hg CAM III B <input checked="" type="checkbox"/>	MassDEP VPH (GC/PID/FID) CAM IV A <input checked="" type="checkbox"/>	8082 PCB CAM V A <input type="checkbox"/>	9014 Total Cyanide/PAC CAM VI A <input type="checkbox"/>	6860 Perchlorate CAM VIII B <input type="checkbox"/>
8270 SVOC CAM II B <input type="checkbox"/>	7010 Metals CAM III C <input type="checkbox"/>	MassDEP VPH (GC/MS) CAM IV C <input type="checkbox"/>	8081 Pesticides CAM V B <input type="checkbox"/>	7196 Hex Cr CAM VI B <input checked="" type="checkbox"/>	MassDEP APH CAM IX A <input type="checkbox"/>
6010 Metals CAM III A <input type="checkbox"/>	6020 Metals CAM III D <input checked="" type="checkbox"/>	MassDEP EPH CAM IV B <input checked="" type="checkbox"/>	8151 Herbicides CAM V C <input type="checkbox"/>	8330 Explosives CAM VIII A <input type="checkbox"/>	TO-15 VOC CAM IX B <input type="checkbox"/>

**Affirmative Responses to Questions A through F are required for "Presumptive Certainty" status**

<b>A</b>	Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature) in the field or laboratory, and prepared/analyzed within method holding times?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>B</b>	Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>C</b>	Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>D</b>	Does the laboratory report comply with all the reporting requirements specified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>E</b>	VPH, EPH, APH, and TO-15 only a. VPH, EPH, and APH Methods only: Was each method conducted without significant modification(s)? (Refer to the individual method(s) for a list of significant modifications). b. APH and TO-15 Methods only: Was the complete analyte list reported for each method?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Yes <input type="checkbox"/> No
<b>F</b>	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No

**Responses to Questions G, H and I below are required for "Presumptive Certainty" status**

<b>G</b>	Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
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**Data User Note:** Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40. 1056 (2)(k) and WSC-07-350.

<b>H</b>	Were all QC performance standards specified in the CAM protocol(s) achieved?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
<b>I</b>	Were results reported for the complete analyte list specified in the selected CAM protocol(s)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <sup>1</sup>

<sup>1</sup>All negative responses must be addressed in an attached laboratory narrative.

*I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, is accurate and complete.*

Signature: 

Position: Chief Operating Officer

Printed Name: Aaron DeWees

Date: 5/1/20



## Sample Integrity Table

Parameter	Method	Matrix	Minimum Volume	Recommended Container(s)	Required Preservation	Holding Time
Volatile Organics	EPA 8260	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
Volatile Organics	EPA 8260	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	14 Days
Semivolatile Organics	EPA 8270	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Semivolatile Organics	EPA 8270	Solid	20g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Organochlorine Pesticides	EPA 8081	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Organochlorine Pesticides	EPA 8081	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
PCBs	EPA 8082	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
PCBs	EPA 8082	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
Herbicides (subcontracted)	EPA 8151	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Herbicides (subcontracted)	EPA 8151	Solid	30g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
MA DEP VPH	MADEP VPH	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP VPH	MADEP VPH	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	28 Days
MA DEP EPH	MADEP EPH	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP EPH	MADEP EPH	Solid	30g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Total Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	180 Days
Dissolved Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	Filter First 1:1 HNO <sub>3</sub> to pH <2	180 Days
Total Metals	EPA 6010	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	180 Days
Total Mercury (may be combined with Total Metals)	EPA 7470	Aqueous	100mL	125mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	28 Days
Total Mercury (may be combined with Total Metals)	EPA 7471	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	28 Days
Chromium, Hexavalent	EPA 7196	Aqueous	100mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> buffer	28 Days
Chromium, Hexavalent (subcontract)	EPA 7196	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	30 Days
Cyanide, Total	EPA 9014	Aqueous	125mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ 1:1 NaOH to pH >8	14 Days
Cyanide, Total	EPA 9014	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days

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**Case Narrative**

**Lab # 52498**

**Sample Receiving and Chain of Custody Discrepancies**

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Samples were received in acceptable condition, between 0 and 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

As noted on the result page, the VPH samples did not meet the 1:1 +/-25% methanol to soil ratio.

**Calibration**

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No exceptions noted.

**Method Blank**

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No exceptions noted.

**Surrogate Recoveries**

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EPH: Sample 52498-003 had recoveries for both extraction surrogates below acceptance criteria and one fractionation surrogate above acceptance criteria. The sample was re-extracted beyond method holding time with similar results. Both extractions are included.

VPH: The surrogate a,a,a-trifluorotoluene was below acceptance criteria in sample 52498-011.

**Laboratory Control Sample Results**

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VOC: The LCSD2001516 did not meet the acceptance criteria for t-butanol (TBA). Since <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

**Matrix Spike/Matrix Spike Duplicate/Duplicate Results**

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Metals: The percent recovery for barium in the matrix spike for sample 52498-002 was 161% the recovery for Zinc was 141%, outside the acceptance criteria of 75-125%. The duplicate of sample 52498-002 had an RPD for barium that was outside the acceptance criteria. Results have been qualified accordingly.

ORP: The duplicate of sample 52498-013 was outside the 10 mV limit. Both results have been included.

**Other**

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VPH: The trap used for VPH analysis is a Tekmar STRATUM Purge Trap 9. The column used for VPH analysis is a Restek Rtx-502.2, 105m, 0.53mmID, and 3um df.

**MassDEP Analytical Protocol Certification Form Questions A through I**

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No explanation is needed for Questions A through I answered in the affirmative.

**Question G:** The CAM protocol reporting limits were not achieved for this project due to dilutions necessary for sample analysis. Box G is "No."

**Question H:** See surrogate section above. Box H is "No."

# GLOSSARY

%R	Percent Recovery
BLK	Blank (Method Blank, Preparation Blank)
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRM	Certified Reference Material (associated with solid Metals samples)
CRMD	Certified Reference Material Duplicate (associated with solid Metals samples)
Dil'n	Dilution
DL	Detection Limit
DUP	Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
LOD	Limit of Detection
LOQ	Limit of Quantitation
MB	Methanol Blank (associated with solid VOC samples)
MLCS	Methanol Laboratory Control Sample (associated with solid VOC samples)
MLCSD	Methanol Laboratory Control Sample Duplicate (associated with solid VOC samples)
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PB	Preparation Blank
QC	Quality Control
RL	Reporting Limit
RPD	Relative Percent Difference
SUR	Surrogate



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- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
MA VPH	MB12628	Unadjusted C5-C8 Aliphatics	<	5.0	ug/g						
		Unadjusted C9-C12 Aliphatics	<	5.0	ug/g						
		methyl t-butyl ether (MTBE)	<	0.10	ug/g						
		benzene	<	0.10	ug/g						
		toluene	<	0.10	ug/g						
		ethylbenzene	<	0.10	ug/g						
		m&p-xylenes	<	0.10	ug/g						
		o-xylene	<	0.10	ug/g						
		naphthalene	<	0.25	ug/g						
		C5-C8 Aliphatics	<	5.0	ug/g						
		C9-C12 Aliphatics	<	5.0	ug/g						
		C9-C10 Aromatics	<	5.0	ug/g						
		2,5-dibromotoluene as Aromatic SUR		113	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		
MA VPH	MLCS12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130		
		Unadjusted C9-C12 Aliphatics		13	ug/g	15	89	70	130		
		methyl t-butyl ether (MTBE)		4.6	ug/g	5	92	70	130		
		benzene		4.8	ug/g	5	97	70	130		
		toluene		4.8	ug/g	5	95	70	130		
		ethylbenzene		4.8	ug/g	5	96	70	130		
		m&p-xylenes		9.7	ug/g	10	97	70	130		
		o-xylene		4.8	ug/g	5	96	70	130		
		naphthalene		5.0	ug/g	5	100	70	130		
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		
		C9-C10 Aromatics		5.1	ug/g	5	102	70	130		
		2,5-dibromotoluene as Aromatic SUR		111	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		86	%			70	130		
MA VPH	MLCSD12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130	0	25
		Unadjusted C9-C12 Aliphatics		14	ug/g	15	90	70	130	2	25
		methyl t-butyl ether (MTBE)		4.7	ug/g	5	93	70	130	1	25
		benzene		4.9	ug/g	5	99	70	130	2	25
		toluene		4.9	ug/g	5	98	70	130	3	25
		ethylbenzene		4.9	ug/g	5	99	70	130	3	25
		m&p-xylenes		10.0	ug/g	10	100	70	130	3	25
		o-xylene		4.9	ug/g	5	99	70	130	2	25
		naphthalene		5.1	ug/g	5	102	70	130	2	25
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C10 Aromatics		5.2	ug/g	5	103	70	130	1	25
		2,5-dibromotoluene as Aromatic SUR		99	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		100	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK12625	naphthalene		<	0.20	ug/g				
		2-methylnaphthalene		<	0.20	ug/g				
		phenanthrene		<	0.20	ug/g				
		acenaphthene		<	0.20	ug/g				
		acenaphthylene		<	0.20	ug/g				
		fluorene		<	0.20	ug/g				
		anthracene		<	0.20	ug/g				
		fluoranthene		<	0.20	ug/g				
		pyrene		<	0.20	ug/g				
		benzo(a)anthracene		<	0.20	ug/g				
		chrysene		<	0.20	ug/g				
		benzo(b)fluoranthene		<	0.20	ug/g				
		benzo(k)fluoranthene		<	0.20	ug/g				
		benzo(a)pyrene		<	0.20	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.20	ug/g				
		dibenzo(a,h)anthracene		<	0.20	ug/g				
		benzo(g,h,i)perylene		<	0.20	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			52	%		40	140	
		o-terphenyl SUR			53	%		40	140	
		2-fluorobiphenyl SUR			67	%		40	140	
		2-bromonaphthalene SUR			65	%		40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCS12625	naphthalene		3.0	ug/g	6	50	40	140	
		2-methylnaphthalene		3.0	ug/g	6	50	40	140	
		phenanthrene		3.6	ug/g	6	60	40	140	
		acenaphthene		3.3	ug/g	6	55	40	140	
		acenaphthylene		3.1	ug/g	6	51	40	140	
		fluorene		3.4	ug/g	6	57	40	140	
		anthracene		3.5	ug/g	6	58	40	140	
		fluoranthene		3.8	ug/g	6	64	40	140	
		pyrene		4.1	ug/g	6	69	40	140	
		benzo(a)anthracene		4.1	ug/g	6	68	40	140	
		chrysene		4.2	ug/g	6	71	40	140	
		benzo(b)fluoranthene		4.0	ug/g	6	67	40	140	
		benzo(k)fluoranthene		3.9	ug/g	6	65	40	140	
		benzo(a)pyrene		4.0	ug/g	6	66	40	140	
		indeno(1,2,3-cd)pyrene		4.2	ug/g	6	71	40	140	
		dibenzo(a,h)anthracene		4.2	ug/g	6	70	40	140	
		benzo(g,h,i)perylene		4.2	ug/g	6	69	40	140	
		Unadjusted C11-C22 Aromatics		68	ug/g	102	66	40	140	
		C9-C18 Aliphatics		20	ug/g	36	57	40	140	
		C19-C36 Aliphatics		56	ug/g	48	116	40	140	
		C11-C22 Aromatics	<	20	ug/g			40	140	
		1-chloro-octadecane SUR		60	%			40	140	
		o-terphenyl SUR		58	%			40	140	
		2-fluorobiphenyl SUR		65	%			40	140	
		2-bromonaphthalene SUR		64	%			40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD12625	naphthalene		2.9	ug/g	6	48	40 140	2	25
		2-methylnaphthalene		2.9	ug/g	6	49	40 140	3	25
		phenanthrene		3.5	ug/g	6	59	40 140	1	25
		acenaphthene		3.3	ug/g	6	55	40 140	0	25
		acenaphthylene		3.1	ug/g	6	51	40 140	0	25
		fluorene		3.4	ug/g	6	57	40 140	1	25
		anthracene		3.4	ug/g	6	57	40 140	1	25
		fluoranthene		3.7	ug/g	6	61	40 140	4	25
		pyrene		4.0	ug/g	6	67	40 140	3	25
		benzo(a)anthracene		4.0	ug/g	6	66	40 140	3	25
		chrysene		4.1	ug/g	6	68	40 140	4	25
		benzo(b)fluoranthene		4.0	ug/g	6	66	40 140	1	25
		benzo(k)fluoranthene		3.8	ug/g	6	63	40 140	3	25
		benzo(a)pyrene		3.9	ug/g	6	64	40 140	3	25
		indeno(1,2,3-cd)pyrene		4.2	ug/g	6	69	40 140	2	25
		dibenzo(a,h)anthracene		4.1	ug/g	6	69	40 140	2	25
		benzo(g,h,i)perylene		4.1	ug/g	6	68	40 140	1	25
		Unadjusted C11-C22 Aromatics		65	ug/g	102	63	40 140	5	25
		C9-C18 Aliphatics	<	20	ug/g	36	52	40 140	8	25
		C19-C36 Aliphatics		54	ug/g	48	113	40 140	3	25
		C11-C22 Aromatics	<	20	ug/g			40 140		25
		1-chloro-octadecane SUR		55	%			40 140		
		o-terphenyl SUR		55	%			40 140		
		2-fluorobiphenyl SUR		68	%			40 140		
		2-bromonaphthalene SUR		66	%			40 140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK12630	naphthalene		<	0.20	ug/g				
		2-methylnaphthalene		<	0.20	ug/g				
		phenanthrene		<	0.20	ug/g				
		acenaphthene		<	0.20	ug/g				
		acenaphthylene		<	0.20	ug/g				
		fluorene		<	0.20	ug/g				
		anthracene		<	0.20	ug/g				
		fluoranthene		<	0.20	ug/g				
		pyrene		<	0.20	ug/g				
		benzo(a)anthracene		<	0.20	ug/g				
		chrysene		<	0.20	ug/g				
		benzo(b)fluoranthene		<	0.20	ug/g				
		benzo(k)fluoranthene		<	0.20	ug/g				
		benzo(a)pyrene		<	0.20	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.20	ug/g				
		dibenzo(a,h)anthracene		<	0.20	ug/g				
		benzo(g,h,i)perylene		<	0.20	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR		52	%			40	140	
		o-terphenyl SUR		59	%			40	140	
		2-fluorobiphenyl SUR		66	%			40	140	
		2-bromonaphthalene SUR		64	%			40	140	



Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCS12630	naphthalene		3.1	ug/g	6	51	40	140	
		2-methylnaphthalene		3.1	ug/g	6	52	40	140	
		phenanthrene		3.6	ug/g	6	60	40	140	
		acenaphthene		3.3	ug/g	6	56	40	140	
		acenaphthylene		3.3	ug/g	6	54	40	140	
		fluorene		3.6	ug/g	6	60	40	140	
		anthracene		3.5	ug/g	6	59	40	140	
		fluoranthene		4.0	ug/g	6	66	40	140	
		pyrene		4.2	ug/g	6	70	40	140	
		benzo(a)anthracene		4.3	ug/g	6	71	40	140	
		chrysene		4.4	ug/g	6	73	40	140	
		benzo(b)fluoranthene		4.4	ug/g	6	73	40	140	
		benzo(k)fluoranthene		3.9	ug/g	6	65	40	140	
		benzo(a)pyrene		4.1	ug/g	6	68	40	140	
		indeno(1,2,3-cd)pyrene		4.4	ug/g	6	73	40	140	
		dibenzo(a,h)anthracene		4.4	ug/g	6	73	40	140	
		benzo(g,h,i)perylene		4.4	ug/g	6	73	40	140	
		Unadjusted C11-C22 Aromatics		65	ug/g	102	64	40	140	
		C9-C18 Aliphatics	<	20	ug/g	36	44	40	140	
		C19-C36 Aliphatics		49	ug/g	48	102	40	140	
		C11-C22 Aromatics	<	20	ug/g			40	140	
		1-chloro-octadecane SUR		52	%			40	140	
		o-terphenyl SUR		57	%			40	140	
		2-fluorobiphenyl SUR		64	%			40	140	
		2-bromonaphthalene SUR		61	%			40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD12630	naphthalene		2.9	ug/g	6	48	40 140	6	25
		2-methylnaphthalene		2.9	ug/g	6	49	40 140	6	25
		phenanthrene		3.5	ug/g	6	59	40 140	2	25
		acenaphthene		3.3	ug/g	6	55	40 140	1	25
		acenaphthylene		3.1	ug/g	6	52	40 140	4	25
		fluorene		3.6	ug/g	6	59	40 140	1	25
		anthracene		3.5	ug/g	6	59	40 140	0	25
		fluoranthene		4.1	ug/g	6	68	40 140	3	25
		pyrene		4.0	ug/g	6	66	40 140	5	25
		benzo(a)anthracene		4.2	ug/g	6	70	40 140	2	25
		chrysene		4.3	ug/g	6	71	40 140	2	25
		benzo(b)fluoranthene		4.1	ug/g	6	68	40 140	8	25
		benzo(k)fluoranthene		4.2	ug/g	6	70	40 140	8	25
		benzo(a)pyrene		4.1	ug/g	6	68	40 140	0	25
		indeno(1,2,3-cd)pyrene		4.2	ug/g	6	71	40 140	4	25
		dibenzo(a,h)anthracene		4.2	ug/g	6	71	40 140	3	25
		benzo(g,h,i)perylene		4.1	ug/g	6	69	40 140	6	25
		Unadjusted C11-C22 Aromatics		67	ug/g	102	66	40 140	3	25
		C9-C18 Aliphatics	<	20	ug/g	36	53	40 140	18	25
		C19-C36 Aliphatics		54	ug/g	48	112	40 140	9	25
		C11-C22 Aromatics	<	20	ug/g			40 140		25
		1-chloro-octadecane SUR		63	%			40 140		
		o-terphenyl SUR		60	%			40 140		
		2-fluorobiphenyl SUR		60	%			40 140		
		2-bromonaphthalene SUR		59	%			40 140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3051A6020A	BLK12626	Silver		<	2.5	ug/g					
		Arsenic		<	2.5	ug/g					
		Barium		<	5.0	ug/g					
		Beryllium		<	0.50	ug/g					
		Cadmium		<	0.50	ug/g					
		Chromium		<	5.0	ug/g					
		Nickel		<	5.0	ug/g					
		Lead		<	2.5	ug/g					
		Antimony		<	0.50	ug/g					
		Selenium		<	5.0	ug/g					
		Thallium		<	0.50	ug/g					
		Vanadium		<	5.0	ug/g					
		Zinc		<	5.0	ug/g					
SW3051A6020A	CRM12626	Silver		47.3	ug/g	53.8		31.9	59.7		
		Arsenic		172	ug/g	219		129	240		
		Barium		634	ug/g	788		509	867		
		Beryllium		204	ug/g	247		160	272		
		Cadmium		155	ug/g	175		111	192		
		Chromium		301	ug/g	375		223	414		
		Nickel		258	ug/g	318		193	358		
		Lead		269	ug/g	321		207	353		
		Antimony		97.6	ug/g	159		15.9	180		
		Selenium		117	ug/g	145		78.5	160		
		Thallium		89.0	ug/g	98.6		53.7	111		
		Vanadium		205	ug/g	267		168	294		
		Zinc		239	ug/g	311		190	352		
SW3051A6020A	CRMD12626	Silver		45.8	ug/g	53.8		31.9	59.7	3	20
		Arsenic		168	ug/g	219		129	240	2	20
		Barium		592	ug/g	788		509	867	7	20
		Beryllium		205	ug/g	247		160	272	1	20
		Cadmium		147	ug/g	175		111	192	6	20
		Chromium		294	ug/g	375		223	414	2	20
		Nickel		254	ug/g	318		193	358	1	20
		Lead		251	ug/g	321		207	353	7	20
		Antimony		92.7	ug/g	159		15.9	180	5	20
		Selenium		106	ug/g	145		78.5	160	10	20
		Thallium		83.4	ug/g	98.6		53.7	111	6	20
		Vanadium		196	ug/g	267		168	294	5	20
		Zinc		230	ug/g	311		190	352	4	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6020A	DUP12626	Silver	52498-002	< 2.6	ug/g					20
		Arsenic	52498-002	4.8	ug/g				19	20
		Barium	52498-002	300	ug/g				39 *	20
		Beryllium	52498-002	< 0.51	ug/g					20
		Cadmium	52498-002	0.53	ug/g				43	20
		Chromium	52498-002	18	ug/g				15	20
		Nickel	52498-002	12	ug/g				3	20
		Lead	52498-002	140	ug/g				13	20
		Antimony	52498-002	0.83	ug/g				9	20
		Selenium	52498-002	< 5.1	ug/g					20
		Thallium	52498-002	< 0.51	ug/g					20
		Vanadium	52498-002	17	ug/g				2	20
		Zinc	52498-002	230	ug/g				17	20
SW3051A6020A	MS12626	Silver	52498-002	120	ug/g	125	97	75	125	
		Arsenic	52498-002	200	ug/g	250	78	75	125	
		Barium	52498-002	610	ug/g	250	161 *	75	125	
		Beryllium	52498-002	240	ug/g	250	96	75	125	
		Cadmium	52498-002	240	ug/g	250	97	75	125	
		Chromium	52498-002	240	ug/g	250	90	75	125	
		Nickel	52498-002	230	ug/g	250	87	75	125	
		Lead	52498-002	400	ug/g	250	110	75	125	
		Antimony	52498-002	240	ug/g	250	94	75	125	
		Selenium	52498-002	200	ug/g	250	78	75	125	
		Vanadium	52498-002	240	ug/g	250	89	75	125	
		Zinc	52498-002	550	ug/g	250	141 *	75	125	
		SW7471B	BLK12622	Mercury		< 0.14	ug/g			
SW7471B	CRM12622	Mercury		0.249	ug/g	0.221		0.0908	0.351	
SW7471B	CRMD12622	Mercury		0.223	ug/g	0.221		0.0908	0.351	11 35
SW7471B	DUP12622	Mercury	52490-001	< 0.14	ug/g					35
SW7471B	MS12622	Mercury	52490-001	0.40	ug/g	0.333	119	80	120	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3060A7196A	CRM2001871	Chromium, Hexavalent		110	ug/g	260		93.3	350	
SW3060A7196A	CRMD2001871	Chromium, Hexavalent		100	ug/g	260		93.3	350	30
SW3060A7196A	DUP2001871	Chromium, Hexavalent	52498-013	< 0.48	ug/g					20
SW3060A7196A	MS2001871	Chromium, Hexavalent	52513-004	< 0.50	ug/g	25.01	1 *	75	125	
SW3060A7196A	MS2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.25	1 *	75	125	
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52513-004	< 0.49	ug/g	24.43	1 *	75	125	25
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.35	1 *	75	125	63
SW3060A7196A	PB2001871	Chromium, Hexavalent		< 0.40	ug/g					
SW9045C	DUP2001823	pH	52498-004	6.0	pH					

## AROMATIC HYDROCARBON BREAKTHROUGH CALCULATION

Method: MADEP EPH 2019 Rev 2.1

	Ics12625 Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.3%	<5.0%	4/8/2020
2-methylnaphthalene	0.2%	<5.0%	4/8/2020

	Icsd12625 Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.3%	<5.0%	4/8/2020
2-methylnaphthalene	0.3%	<5.0%	4/8/2020

## AROMATIC HYDROCARBON BREAKTHROUGH CALCULATION

Method: MADEP EPH 2019 Rev 2.1

	Ics12630	Acceptance	Date of Analysis
	Aliphatic Breakthrough	Criteria	
	(%)		
naphthalene	0.6%	<5.0%	4/9/2020
2-methylnaphthalene	0.2%	<5.0%	4/9/2020

	Icsd12630	Acceptance	Date of Analysis
	Aliphatic Breakthrough	Criteria	
	(%)		
naphthalene	0.4%	<5.0%	4/9/2020
2-methylnaphthalene	0.2%	<5.0%	4/9/2020

**Absolute Resource**  
associates



124 Heritage Avenue #16  
Portsmouth, NH 03801  
603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
AND ANALYSIS REQUEST**

52498

**ANALYSIS REQUEST**

Company Name: **WESTON & SAMPSON**  
Company Address: **55 WALKERS BROOK DR READING, MA**  
Report To: **SARAH DESTEFANO & JILL MURPHY**  
Phone #: **1 800 SAMPSON**  
Invoice to: **SARAH DESTEFANO**  
Email: **DESTEFANOS@WSEINC.COM**  
PO #: **FARMINGHAM BROWNFIELDS**

Project Name: **CEDAR WOODS**  
Project #: **2180311**  
Project Location: **NH (MA) ME VT**  
Accreditation Required? N/Y: **Y**  
Protocol: **RCRA MCP SDWA NHDES NPDES DOD**  
Reporting Limits: **EPA DW Other (S-1)**  
Quote #: **N/A**  
 NH Reimbursement Pricing

<input type="checkbox"/> VOC 8260	<input type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MADEP	<input type="checkbox"/> VOC 624.1	<input type="checkbox"/> VOC BTEX MIBE, only	<input type="checkbox"/> VOC 8021VT	<input checked="" type="checkbox"/> VPH MADEP	<input type="checkbox"/> GRO 8015	<input checked="" type="checkbox"/> 1,4-Dioxane *	<input type="checkbox"/> VOC 524.2	<input type="checkbox"/> VOC 524.2 NH List	<input type="checkbox"/> Gases-List:	<input type="checkbox"/> TPH	<input type="checkbox"/> DR0 8015	<input checked="" type="checkbox"/> EPH MADEP	<input type="checkbox"/> TPH Fingerprint	<input type="checkbox"/> 8270PAH	<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625.1	<input type="checkbox"/> EDB	<input type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608.3 Pest/PCB	<input type="checkbox"/> O&G 1664	<input type="checkbox"/> Mineral O&G 1664	<input type="checkbox"/> pH	<input type="checkbox"/> BOD	<input type="checkbox"/> Conductivity	<input type="checkbox"/> Turbidity	<input type="checkbox"/> Apparent Color	<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TS	<input type="checkbox"/> TVS	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Acidity	<input type="checkbox"/> RCRA Metals	<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input type="checkbox"/> Hardness	<input checked="" type="checkbox"/> Total Metals-list: <b>MCP 14 METALS</b>	<input type="checkbox"/> Dissolved Metals-list:	<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD	<input type="checkbox"/> TKN	<input type="checkbox"/> TN	<input type="checkbox"/> TON	<input type="checkbox"/> TOC	<input type="checkbox"/> Ferrous Iron	<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Bacteria P/A	<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> Enterococci	<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide	<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Ortho P	<input type="checkbox"/> Phenols	<input type="checkbox"/> Nitrate	<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Sulfate	<input type="checkbox"/> Bromide	<input type="checkbox"/> Fluoride	<input type="checkbox"/> Corrosivity	<input type="checkbox"/> Ignitibility/FP	<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC	<input type="checkbox"/> TCLP Pesticide	<input type="checkbox"/> Subcontract: <input type="checkbox"/> Grain Size <input type="checkbox"/> Herbicides <input type="checkbox"/> Asbestos <input type="checkbox"/> PFAS	<b>HEXAVALENT CHROMIUM (Cr VI)</b>	<input type="checkbox"/> Grab (G) or Composite (C)
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Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method				Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME
52498-01	SB-106(0-1')			X					X	4/6/20	0920	
-02	SB-106(0-3')			X					X		0940	
-03	SB-106(5-8')			X					X		1000	
-04	SB-106(3-4')			X					X		0950	
-05	SB-105(0-3')			X					X		1040	
-06	SB-105(5-8')			X					X		1100	
-07	SB-102(0-3')			X					X		1145	
-08	SB-102(0-1')			X					X		1130	
-09	SB-102(11-14')			X					X		1140	
-10	SB-101(0-3')			X					X		1255	
-11	SB-101(9-12')			X					X		1305	

**TAT REQUESTED**  
Priority (24 hr)\*   
Expedited (48 hr)\*   
Standard (10 Business Days)   
\*Date Needed: **5-DAY**

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
\*VPH W/ TARGET VOCs

REPORTING INSTRUCTIONS:  PDF (e-mail address) **murphy.j@wseinc.com**  
 HARD COPY REQUIRED  EDD

RECEIVED ON ICE:  YES  NO  
TEMPERATURE: **0** °C

**CUSTODY RECORD**  
QSD-01 Revision 11/06/19

Relinquished by Sampler:	Date: <b>4/6/20</b>	Time: <b>1340</b>	Received by:	Date: <b>4-6</b>	Time: <b>1:41</b>
Relinquished by:	Date: <b>4-6</b>	Time: <b>15:43</b>	Received by:	Date:	Time:
Relinquished by:	Date:	Time:	Received by Laboratory:	Date: <b>4/6/20</b>	Time: <b>15:43</b>



**Absolute Resource**  
associates



124 Heritage Avenue #16  
Portsmouth, NH 03801  
603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
AND ANALYSIS REQUEST**

52498

**ANALYSIS REQUEST**

Company Name: \_\_\_\_\_ Project Name: \_\_\_\_\_  
 Company Address: **SEE PAGE 1 OF 2** Project #: \_\_\_\_\_  
 Report To: \_\_\_\_\_ Accreditation Required? N/Y: \_\_\_\_\_  
 Phone #: \_\_\_\_\_ Protocol: RCRA SDWA NPDES  
 Reporting QAPP GW-1 S-1  
 Limits: EPA DW Other \_\_\_\_\_  
 Invoice to: \_\_\_\_\_ Quote # \_\_\_\_\_  
 Email: \_\_\_\_\_  
 PO #: \_\_\_\_\_  NH Reimbursement Pricing

VOC 8260  VOC 8260 NHDES  VOC 8260 MADEP  
 VOC 624.1  VOC BTEX MIBE, only  VOC 8021VT  
 VPH MADEP  GRO 8015  1,4-Dioxane \*  
 VOC 524.2  VOC 524.2 NH List  Gases-List:  
 TPH  DRO 8015  VEPH MADEP  TPH Fingerprint  
 8270PAH  8270ABN  625.1  EDB  
 8082 PCB  8081 Pesticides  608.3 Pest/PCB  
 O&G 1664  Mineral O&G 1664  
 pH  80D  Conductivity  Turbidity  Apparent Color  
 TSS  TDS  TS  TVS  Alkalinity  Acidity  
 RCRA Metals  Priority Pollutant Metals  TAL Metals  Hardness  
 Total Metals-list: **MCP 14 METALS**  
 Dissolved Metals-list:  
 Ammonia  COD  TKN  TN  TOC  Ferrrous Iron  
 T-Phosphorus  Bacteria P/A  Bacteria MPN  Enterococci  
 Cyanide  Sulfide  Nitrate + Nitrite  Ortho P  Phenols  
 Nitrate  Nitrite  Chloride  Sulfate  Bromide  Fluoride  
 Corrosivity  Ignitibility/FP  
 TCLP Metals  TCLP VOC  TCLP SVOC  TCLP Pesticide  
 Subcontract:  Grain Size  Herbicides  Asbestos  PFAS  
**XX HEXAVALENT CHROMIUM (CR VI)**  
 Grab (G) or Composite (C)

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
52498-12	SB-103(0-3')			X					X	4/6/20	1220	X	
-13	SB-103(3-5')			X					X	↓	1220	X	
-14	SB-103(7-10')			X					X		1225	X	
-15	TREP BLANK			X					X	4/6/20		X	

**TAT REQUESTED**  
 Priority (24 hr)\*   
 Expedited (48 hr)\*   
 Standard (10 Business Days)   
 \*Date Needed: **5-DAY**

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
 \*VPH w/TARGET VOCs

**REPORTING INSTRUCTIONS**  PDF (e-mail address) \_\_\_\_\_  
 HARD COPY REQUIRED  EDD

RECEIVED ON ICE  YES  NO  
 TEMPERATURE **0** °C

**CUSTODY RECORD**  
 QSD-01 Revision 11/06/19

Relinquished by Sampler: _____	Date: 4/6/20	Time: 1340	Received by: _____	Date: 4-6	Time: 1-40
Relinquished by: _____	Date: 4-6	Time: 1543	Received by: _____	Date: 4/6/20	Time: 1543
Relinquished by: _____	Date: _____	Time: _____	Received by Laboratory: _____	Date: _____	Time: _____

# Laboratory Report



**Absolute Resource** *associates*

124 Heritage Avenue Portsmouth NH 03801

Jill Murphy  
Weston & Sampson  
55 Walkers Brook Drive  
Reading, MA 01867

PO Number: Framingham Brownfields  
Job ID: 52513  
Date Received: 4/7/20

Project: Cedar Woods 21

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below. The reported results apply to the sample(s) in the condition as received at the time the laboratory took custody. This report shall not be reproduced except in full and with approval from the laboratory. The liability of ARA is limited to the cost of the requested analyses, unless otherwise agreed upon in writing.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,  
Absolute Resource Associates

A handwritten signature in black ink, appearing to read 'A. DeWees', written in a cursive style.

Aaron DeWees  
Chief Operating Officer

Date of Approval: 5/4/2020  
Total number of pages: 69

## Absolute Resource Associates Certifications

New Hampshire 1732  
Maine NH902

Massachusetts M-NH902

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-104 (0-3')	Solid	4/7/2020 8:45	52513-001	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-104 (8-10')	Solid	4/7/2020 8:55	52513-002	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-107 (0-3')	Solid	4/7/2020 10:00	52513-004	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Field Specified Laboratory Duplicate Field Specified Matrix Spike Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
SB-107 (10-11')	Solid	4/7/2020 10:20	52513-005	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-108 (0-3')	Solid	4/7/2020 10:50	52513-007	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-108 (6-9')	Solid	4/7/2020 11:00	52513-008	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VOCs in solids by 8260 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-109 (0-3')	Solid	4/7/2020 11:40	52513-010	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-109 (5-8')	Solid	4/7/2020 11:45	52513-011	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-110 (0-3')	Solid	4/7/2020 12:50	52513-013	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-110 (5-8')	Solid	4/7/2020 12:55	52513-014	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-112 (0-3')	Solid	4/7/2020 13:20	52513-016	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-112 (5-8')	Solid	4/7/2020 13:25	52513-017	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020



## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
DUP-1	Solid	4/7/2020 0:00	52513-018	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
Trip Blank	Solid	4/7/2020 0:00	52513-019	VOCs in solids by 8260

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-008

Sample ID: SB-108 (6-9')

Matrix: Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 11:00

Parameter	Result	Reporting		Instr Dil'n		Prep		Analysis		Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
dichlorodifluoromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
chloromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
vinyl chloride	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
bromomethane	< 0.27	0.27	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
chloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
trichlorofluoromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
diethyl ether	< 0.54	0.54	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
acetone	< 2.7	2.7	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1-dichloroethene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
methylene chloride	< 0.27	0.27	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
carbon disulfide	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
methyl t-butyl ether (MTBE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
trans-1,2-dichloroethene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
isopropyl ether (DIPE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
ethyl t-butyl ether (ETBE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1-dichloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
t-butanol (TBA)	< 2.7	2.7	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
2-butanone (MEK)	< 0.32	0.32	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
2,2-dichloropropane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
cis-1,2-dichloroethene	<b>0.20</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
chloroform	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
bromochloromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
tetrahydrofuran (THF)	< 0.54	0.54	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1,1-trichloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1-dichloropropene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
t-amyl-methyl ether (TAME)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
carbon tetrachloride	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2-dichloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
benzene	<b>0.19</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
trichloroethene	<b>0.36</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2-dichloropropane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
bromodichloromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,4-dioxane	< 2.7	2.7	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
dibromomethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
4-methyl-2-pentanone (MIBK)	< 0.49	0.49	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
cis-1,3-dichloropropene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
toluene	<b>0.46</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
trans-1,3-dichloropropene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
2-hexanone	< 0.54	0.54	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1,2-trichloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,3-dichloropropane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-008

Sample ID: SB-108 (6-9')

Matrix: Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 11:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis			Reference	
	Result	Limit				Analyst	Date	Batch		Date
tetrachloroethene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
dibromochloromethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2-dibromoethane (EDB)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
chlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1,1,2-tetrachloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
ethylbenzene	<b>0.59</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
m&p-xylenes	<b>2.9</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
o-xylene	<b>0.53</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
styrene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
bromoform	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
isopropylbenzene	<b>0.34</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,1,2,2-tetrachloroethane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2,3-trichloropropane	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
n-propylbenzene	<b>0.93</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
bromobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,3,5-trimethylbenzene	<b>2.3</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
2-chlorotoluene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
4-chlorotoluene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
tert-butylbenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2,4-trimethylbenzene	<b>6.9</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
sec-butylbenzene	<b>0.26</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,3-dichlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
4-isopropyltoluene	<b>0.30</b>	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,4-dichlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2-dichlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
n-butylbenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2-dibromo-3-chloropropane (DBCP)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2,4-trichlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
hexachlorobutadiene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
naphthalene	<b>1.8</b>	0.27	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
1,2,3-trichlorobenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
<b>Surrogate Recovery</b>		<b>Limits</b>								
dibromofluoromethane SUR	<b>81</b>	78-114	%	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
toluene-D8 SUR	<b>98</b>	88-110	%	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
4-bromofluorobenzene SUR	<b>116 *</b>	86-115	%	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D
a,a,a-trifluorotoluene SUR	<b>89</b>	70-130	%	1	LMM	4/8/20	12627	4/10/20	9:36	SW5035A8260D

\* The surrogate showed recovery outside the acceptance limits as a result of hydrocarbons present in the sample.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-019

Sample ID: Trip Blank

Matrix: Solid

Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 0:00

Parameter	Result	Reporting		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
dichlorodifluoromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
chloromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
vinyl chloride	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
bromomethane	< 0.25	0.25	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
chloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
trichlorofluoromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
diethyl ether	< 0.50	0.50	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
acetone	< 2.5	2.5	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1-dichloroethene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
methylene chloride	< 0.25	0.25	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
carbon disulfide	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
methyl t-butyl ether (MTBE)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
trans-1,2-dichloroethene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
isopropyl ether (DIPE)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
ethyl t-butyl ether (ETBE)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1-dichloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
t-butanol (TBA)	< 2.5	2.5	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
2-butanone (MEK)	< 0.30	0.30	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
2,2-dichloropropane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
cis-1,2-dichloroethene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
chloroform	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
bromochloromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
tetrahydrofuran (THF)	< 0.50	0.50	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1,1-trichloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1-dichloropropene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
t-amyl-methyl ether (TAME)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
carbon tetrachloride	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2-dichloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
benzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
trichloroethene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2-dichloropropane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
bromodichloromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,4-dioxane	< 2.5	2.5	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
dibromomethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
4-methyl-2-pentanone (MIBK)	< 0.45	0.45	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
cis-1,3-dichloropropene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
toluene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
trans-1,3-dichloropropene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
2-hexanone	< 0.50	0.50	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1,2-trichloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,3-dichloropropane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-019

Sample ID: Trip Blank

Matrix: Solid

Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 0:00

Parameter	Result	Reporting		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
tetrachloroethene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
dibromochloromethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2-dibromoethane (EDB)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
chlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1,1,2-tetrachloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
ethylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
m&p-xylenes	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
o-xylene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
styrene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
bromoform	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
isopropylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,1,2,2-tetrachloroethane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2,3-trichloropropane	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
n-propylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
bromobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,3,5-trimethylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
2-chlorotoluene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
4-chlorotoluene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
tert-butylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2,4-trimethylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
sec-butylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,3-dichlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
4-isopropyltoluene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,4-dichlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2-dichlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
n-butylbenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2-dibromo-3-chloropropane (DBCP)	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2,4-trichlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
hexachlorobutadiene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
naphthalene	< 0.25	0.25	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
1,2,3-trichlorobenzene	< 0.10	0.10	ug/g	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
<b>Surrogate Recovery</b>		<b>Limits</b>								
dibromofluoromethane SUR	<b>98</b>	78-114	%	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
toluene-D8 SUR	<b>105</b>	88-110	%	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
4-bromofluorobenzene SUR	<b>110</b>	86-115	%	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D
a,a,a-trifluorotoluene SUR	<b>110</b>	70-130	%	1	LMM	4/8/20	12627	4/9/20	12:48	SW5035A8260D

**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-002

**Sample ID:** SB-104 (8-10')

**Matrix:** Solid Percent Dry: 83.1% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.64 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

**Sampled:** 4/7/20 8:55

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Analyst	Date	Batch	Date	Time
Unadjusted C5-C8 Aliphatics	< 4.9	4.9	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 4.9	4.9	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
benzene	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
toluene	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
ethylbenzene	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
m&p-xylenes	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
o-xylene	< 0.097	0.097	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
naphthalene	< 0.24	0.24	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
C5-C8 Aliphatics	< 4.9	4.9	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
C9-C12 Aliphatics	< 4.9	4.9	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
C9-C10 Aromatics	< 4.9	4.9	ug/g	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>106</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>107</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>79</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-004

Sample ID: SB-107 (0-3')

Matrix: Solid Percent Dry: 79% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.65 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	7.6	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
benzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
toluene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
ethylbenzene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
m&p-xylenes	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
o-xylene	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
naphthalene	< 0.27	0.27	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
C5-C8 Aliphatics	7.6	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
C9-C12 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
C9-C10 Aromatics	< 5.4	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	106	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	106	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH
a,a,a-trifluorotoluene SUR	32 *	70-130	%	1	LMM	4/8/20	12628	4/9/20	15:49	MA VPH

\* This surrogate showed recovery outside the acceptance limits.

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-005

Sample ID: SB-107 (10-11')

Matrix: Solid Percent Dry: 84.5% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.64 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 10:20

Parameter	Reporting		Instr Dil'n	Prep	Analysis			Reference		
	Result	Limit			Units	Factor	Analyst		Date	Batch
Unadjusted C5-C8 Aliphatics	26	4.7	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
Unadjusted C9-C12 Aliphatics	62	4.7	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.094	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
benzene	< 0.094	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
toluene	< 0.094	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
ethylbenzene	0.48	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
m&p-xylenes	1.6	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
o-xylene	< 0.094	0.094	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
naphthalene	1.3	0.23	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
C5-C8 Aliphatics	26	4.7	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
C9-C12 Aliphatics	22	4.7	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
C9-C10 Aromatics	39	4.7	ug/g	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	121	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	136 *	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH
a,a,a-trifluorotoluene SUR	115	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:19	MA VPH

\* The surrogate showed recovery outside the acceptance limits as a result of hydrocarbons present in the sample.

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.



Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-008

Sample ID: SB-108 (6-9')

Matrix: Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 11:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	31	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
Unadjusted C9-C12 Aliphatics	95	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.11	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
benzene	0.30	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
toluene	0.42	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
ethylbenzene	0.49	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
m&p-xylenes	2.7	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
o-xylene	0.45	0.11	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
naphthalene	1.9	0.27	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
C5-C8 Aliphatics	30	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
C9-C12 Aliphatics	24	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
C9-C10 Aromatics	67	5.4	ug/g	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	98	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	98	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH
a,a,a-trifluorotoluene SUR	96	70-130	%	1	LMM	4/8/20	12628	4/9/20	16:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-011

**Sample ID:** SB-109 (5-8')

**Matrix:** Solid      Percent Dry: 23.7% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.63 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

**Sampled:** 4/7/20 11:45

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Analyst	Date	Batch	Date	Time
Unadjusted C5-C8 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
benzene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
toluene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
ethylbenzene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
m&p-xylenes	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
o-xylene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
naphthalene	< 1.5	1.5	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
C5-C8 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
C9-C12 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
C9-C10 Aromatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>100</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>98</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>75</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-014

Sample ID: SB-110 (5-8')

Matrix: Solid Percent Dry: 17.7% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.65 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 12:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 42	42	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 42	42	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
benzene	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
toluene	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
ethylbenzene	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
m&p-xylenes	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
o-xylene	< 0.83	0.83	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
naphthalene	< 2.1	2.1	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
C5-C8 Aliphatics	< 42	42	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
C9-C12 Aliphatics	< 42	42	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
C9-C10 Aromatics	< 42	42	ug/g	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>102</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>103</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>158 *</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	17:49	MA VPH

**\* This surrogate showed recovery outside the acceptance limits. This is likely a result of high moisture content in the sample.**

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-017

**Sample ID:** SB-112 (5-8')

**Matrix:** Solid      Percent Dry: 52.4% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.61 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

**Sampled:** 4/7/20    13:25

Parameter	Reporting		Instr Dil'n	Prep	Analysis			Reference		
	Result	Limit			Units	Factor	Analyst		Date	Batch
Unadjusted C5-C8 Aliphatics	< 10	10	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 10	10	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
benzene	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
toluene	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
ethylbenzene	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
m&p-xylenes	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
o-xylene	< 0.21	0.21	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
naphthalene	< 0.52	0.52	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
C5-C8 Aliphatics	< 10	10	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
C9-C12 Aliphatics	< 10	10	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
C9-C10 Aromatics	< 10	10	ug/g	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>105</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>106</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>75</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-018

Sample ID: DUP-1

Matrix: Solid Percent Dry: 23.7% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.63 mL MeOH/g soil.

Received on ice at 2°C, in satisfactory condition.

Sampled: 4/7/20 0:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis			Reference	
	Result	Limit				Analyst	Date	Batch		Date
Unadjusted C5-C8 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
benzene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
toluene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
ethylbenzene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
m&p-xylenes	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
o-xylene	< 0.59	0.59	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
naphthalene	< 1.5	1.5	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
C5-C8 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
C9-C12 Aliphatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
C9-C10 Aromatics	< 29	29	ug/g	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>107</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>105</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>71</b>	70-130	%	1	LMM	4/8/20	12628	4/9/20	18:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-001

Sample ID: SB-104 (0-3')

Matrix: Solid

Percent Dry: 89.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 8:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
2-methylnaphthalene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
phenanthrene	<b>0.39</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
acenaphthene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
acenaphthylene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
fluorene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
anthracene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
fluoranthene	<b>0.80</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
pyrene	<b>0.85</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
benzo(a)anthracene	<b>0.40</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
chrysene	<b>0.56</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
benzo(b)fluoranthene	<b>0.46</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
benzo(k)fluoranthene	<b>0.44</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
benzo(a)pyrene	<b>0.44</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
indeno(1,2,3-cd)pyrene	<b>0.30</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
dibenzo(a,h)anthracene	< 0.20	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
benzo(g,h,i)perylene	<b>0.35</b>	0.20	ug/g	1	CL	4/8/20	12630	4/9/20	18:13	MA EPH
Unadjusted C11-C22 Aromatics	<b>31</b>	20	ug/g	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
C9-C18 Aliphatics	< 20	20	ug/g	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
C19-C36 Aliphatics	<b>34</b>	20	ug/g	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
C11-C22 Aromatics	<b>25</b>	20	ug/g	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>55</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
o-terphenyl SUR	<b>61</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
2-fluorobiphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH
2-bromonaphthalene SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	12:48	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-002

Sample ID: SB-104 (8-10')

Matrix: Solid

Percent Dry: 83.1% Results expressed on a dry weight basis.

Sampled: 4/7/20 8:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
2-methylnaphthalene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
phenanthrene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
acenaphthene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
acenaphthylene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
fluorene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
anthracene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
fluoranthene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
pyrene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
benzo(a)anthracene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
chrysene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
benzo(b)fluoranthene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
benzo(k)fluoranthene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
benzo(a)pyrene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
indeno(1,2,3-cd)pyrene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
dibenzo(a,h)anthracene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
benzo(g,h,i)perylene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	19:44	MA EPH
Unadjusted C11-C22 Aromatics	< 22	22	ug/g	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
C9-C18 Aliphatics	< 22	22	ug/g	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
C19-C36 Aliphatics	<b>100</b>	22	ug/g	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
C11-C22 Aromatics	< 22	22	ug/g	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
o-terphenyl SUR	<b>71</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
2-fluorobiphenyl SUR	<b>70</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH
2-bromonaphthalene SUR	<b>69</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:55	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-004

Sample ID: SB-107 (0-3')

Matrix: Solid Percent Dry: 79% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
2-methylnaphthalene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
phenanthrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
acenaphthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
acenaphthylene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
fluorene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
fluoranthene	<b>0.25</b>	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
benzo(a)anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
chrysene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
benzo(b)fluoranthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
benzo(k)fluoranthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
benzo(a)pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
indeno(1,2,3-cd)pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
dibenzo(a,h)anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
benzo(g,h,i)perylene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	14:41	MA EPH
Unadjusted C11-C22 Aromatics	< 23	23	ug/g	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
C9-C18 Aliphatics	< 23	23	ug/g	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
C19-C36 Aliphatics	<b>36</b>	23	ug/g	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
C11-C22 Aromatics	< 23	23	ug/g	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>55</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
o-terphenyl SUR	<b>54</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
2-fluorobiphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH
2-bromonaphthalene SUR	<b>59</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:22	MA EPH



Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-005

Sample ID: SB-107 (10-11')

Matrix: Solid Percent Dry: 84.5% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
2-methylnaphthalene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
phenanthrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
acenaphthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
acenaphthylene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
fluorene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
fluoranthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
benzo(a)anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
chrysene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
benzo(b)fluoranthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
benzo(k)fluoranthene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
benzo(a)pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
indeno(1,2,3-cd)pyrene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
dibenzo(a,h)anthracene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
benzo(g,h,i)perylene	< 0.23	0.23	ug/g	1	CL	4/8/20	12630	4/9/20	18:44	MA EPH
Unadjusted C11-C22 Aromatics	<b>77</b>	23	ug/g	1	DBV	4/8/20	12630	4/9/20	21:21	MA EPH
C9-C18 Aliphatics	<b>34</b>	23	ug/g	1	DBV	4/8/20	12630	4/14/20	12:58	MA EPH
C19-C36 Aliphatics	<b>410</b>	23	ug/g	1	DBV	4/8/20	12630	4/14/20	12:58	MA EPH
C11-C22 Aromatics	<b>76</b>	23	ug/g	1	DBV	4/8/20	12630	4/9/20	21:21	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>60</b>	40-140	%	1	DBV	4/8/20	12630	4/14/20	12:58	MA EPH
o-terphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:21	MA EPH
2-fluorobiphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:21	MA EPH
2-bromonaphthalene SUR	<b>64</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	21:21	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-007

Sample ID: SB-108 (0-3')

Matrix: Solid Percent Dry: 73.6% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:50

Parameter	Result	Reporting		Instr Dil'n		Prep		Analysis		
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
2-methylnaphthalene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
phenanthrene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
acenaphthene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
acenaphthylene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
fluorene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
anthracene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
fluoranthene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
pyrene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
benzo(a)anthracene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
chrysene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
benzo(b)fluoranthene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
benzo(k)fluoranthene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
benzo(a)pyrene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
indeno(1,2,3-cd)pyrene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
dibenzo(a,h)anthracene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
benzo(g,h,i)perylene	< 0.26	0.26	ug/g	1	CL	4/8/20	12630	4/9/20	15:11	MA EPH
Unadjusted C11-C22 Aromatics	< 26	26	ug/g	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
C9-C18 Aliphatics	< 26	26	ug/g	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
C19-C36 Aliphatics	<b>39</b>	26	ug/g	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
C11-C22 Aromatics	< 26	26	ug/g	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>60</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
o-terphenyl SUR	<b>57</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
2-fluorobiphenyl SUR	<b>60</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH
2-bromonaphthalene SUR	<b>58</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	13:56	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-008

Sample ID: SB-108 (6-9')

Matrix: Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	0.50	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
2-methylnaphthalene	0.57	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
phenanthrene	2.1	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
acenaphthene	< 0.24	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
acenaphthylene	< 0.24	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
fluorene	0.37	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
anthracene	0.46	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
fluoranthene	1.9	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
pyrene	2.1	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
benzo(a)anthracene	0.79	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
chrysene	0.87	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
benzo(b)fluoranthene	0.64	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
benzo(k)fluoranthene	0.57	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
benzo(a)pyrene	0.70	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
indeno(1,2,3-cd)pyrene	0.36	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
dibenzo(a,h)anthracene	< 0.24	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
benzo(g,h,i)perylene	0.51	0.24	ug/g	1	CL	4/8/20	12630	4/9/20	20:15	MA EPH
Unadjusted C11-C22 Aromatics	360	24	ug/g	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
C9-C18 Aliphatics	92	24	ug/g	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
C19-C36 Aliphatics	1900	24	ug/g	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
C11-C22 Aromatics	350	24	ug/g	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	58	40-140	%	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
o-terphenyl SUR	62	40-140	%	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
2-fluorobiphenyl SUR	68	40-140	%	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH
2-bromonaphthalene SUR	72	40-140	%	1	DBV	4/8/20	12630	4/9/20	20:46	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-010

Sample ID: SB-109 (0-3')

Matrix: Solid Percent Dry: 70.9% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
2-methylnaphthalene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
phenanthrene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
acenaphthene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
acenaphthylene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
fluorene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
anthracene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
fluoranthene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
pyrene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
benzo(a)anthracene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
chrysene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
benzo(b)fluoranthene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
benzo(k)fluoranthene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
benzo(a)pyrene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
indeno(1,2,3-cd)pyrene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
dibenzo(a,h)anthracene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
benzo(g,h,i)perylene	< 0.28	0.28	ug/g	1	CL	4/8/20	12630	4/9/20	15:41	MA EPH
Unadjusted C11-C22 Aromatics	< 28	28	ug/g	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
C9-C18 Aliphatics	< 28	28	ug/g	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
C19-C36 Aliphatics	<b>42</b>	28	ug/g	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
C11-C22 Aromatics	< 28	28	ug/g	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>56</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
o-terphenyl SUR	<b>61</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
2-fluorobiphenyl SUR	<b>68</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH
2-bromonaphthalene SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	14:31	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-011

Sample ID: SB-109 (5-8')

Matrix: Solid Percent Dry: 23.7% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
2-methylnaphthalene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
phenanthrene	<b>0.80</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
acenaphthene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
acenaphthylene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
fluorene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
anthracene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
fluoranthene	<b>1.5</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
pyrene	<b>1.7</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
benzo(a)anthracene	<b>0.78</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
chrysene	<b>0.96</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
benzo(b)fluoranthene	<b>0.89</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
benzo(k)fluoranthene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
benzo(a)pyrene	<b>0.91</b>	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
indeno(1,2,3-cd)pyrene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
dibenzo(a,h)anthracene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
benzo(g,h,i)perylene	< 0.78	0.78	ug/g	1	CL	4/8/20	12630	4/9/20	16:12	MA EPH
Unadjusted C11-C22 Aromatics	<b>110</b>	78	ug/g	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
C9-C18 Aliphatics	< 78	78	ug/g	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
C19-C36 Aliphatics	<b>290</b>	78	ug/g	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
C11-C22 Aromatics	<b>100</b>	78	ug/g	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>62</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
o-terphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
2-fluorobiphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH
2-bromonaphthalene SUR	<b>60</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:05	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-013

Sample ID: SB-110 (0-3')

Matrix: Solid Percent Dry: 88.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 12:50

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
2-methylnaphthalene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
phenanthrene	1.1	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
acenaphthene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
acenaphthylene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
fluorene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
anthracene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
fluoranthene	1.3	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
pyrene	1.4	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
benzo(a)anthracene	0.59	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
chrysene	0.74	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
benzo(b)fluoranthene	0.57	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
benzo(k)fluoranthene	0.53	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
benzo(a)pyrene	0.63	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
indeno(1,2,3-cd)pyrene	0.36	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
dibenzo(a,h)anthracene	< 0.22	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
benzo(g,h,i)perylene	0.42	0.22	ug/g	1	CL	4/8/20	12630	4/9/20	16:42	MA EPH
Unadjusted C11-C22 Aromatics	62	22	ug/g	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
C9-C18 Aliphatics	< 22	22	ug/g	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
C19-C36 Aliphatics	68	22	ug/g	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
C11-C22 Aromatics	53	22	ug/g	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	45	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
o-terphenyl SUR	54	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
2-fluorobiphenyl SUR	73	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH
2-bromonaphthalene SUR	72	40-140	%	1	DBV	4/8/20	12630	4/9/20	15:39	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-014

Sample ID: SB-110 (5-8')

Matrix: Solid

Percent Dry: 17.7% Results expressed on a dry weight basis.

Sampled: 4/7/20 12:55

Parameter	Result	Reporting		Instr Dil'n		Prep		Analysis			Reference
		Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
naphthalene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
2-methylnaphthalene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
phenanthrene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
acenaphthene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
acenaphthylene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
fluorene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
anthracene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
fluoranthene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
pyrene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
benzo(a)anthracene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
chrysene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
benzo(b)fluoranthene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
benzo(k)fluoranthene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
benzo(a)pyrene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
indeno(1,2,3-cd)pyrene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
dibenzo(a,h)anthracene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
benzo(g,h,i)perylene	< 1.1	1.1	ug/g	1	CL	4/8/20	12630	4/9/20	14:10	MA EPH	
Unadjusted C11-C22 Aromatics	< 110	110	ug/g	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
C9-C18 Aliphatics	< 110	110	ug/g	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
C19-C36 Aliphatics	<b>140</b>	110	ug/g	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
C11-C22 Aromatics	< 110	110	ug/g	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
<b>Surrogate Recovery</b>		<b>Limits</b>									
1-chloro-octadecane SUR	<b>44</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
o-terphenyl SUR	<b>47</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
2-fluorobiphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	
2-bromonaphthalene SUR	<b>62</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:13	MA EPH	

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-016

Sample ID: SB-112 (0-3')

Matrix: Solid Percent Dry: 90.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 13:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
2-methylnaphthalene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
phenanthrene	<b>0.41</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
acenaphthene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
acenaphthylene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
fluorene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
anthracene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
fluoranthene	<b>0.80</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
pyrene	<b>0.73</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
benzo(a)anthracene	<b>0.39</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
chrysene	<b>0.42</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
benzo(b)fluoranthene	<b>0.37</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
benzo(k)fluoranthene	<b>0.38</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
benzo(a)pyrene	<b>0.40</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
indeno(1,2,3-cd)pyrene	<b>0.21</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
dibenzo(a,h)anthracene	< 0.21	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
benzo(g,h,i)perylene	<b>0.25</b>	0.21	ug/g	1	CL	4/8/20	12630	4/9/20	17:12	MA EPH
Unadjusted C11-C22 Aromatics	<b>26</b>	21	ug/g	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
C9-C18 Aliphatics	< 21	21	ug/g	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
C19-C36 Aliphatics	<b>50</b>	21	ug/g	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
C11-C22 Aromatics	<b>22</b>	21	ug/g	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>62</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
o-terphenyl SUR	<b>68</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
2-fluorobiphenyl SUR	<b>66</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH
2-bromonaphthalene SUR	<b>63</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	16:47	MA EPH



Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-017

Sample ID: SB-112 (5-8')

Matrix: Solid

Percent Dry: 52.4% Results expressed on a dry weight basis.

Sampled: 4/7/20 13:25

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.38	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
2-methylnaphthalene	< 0.38	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
phenanthrene	<b>4.7</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
acenaphthene	< 0.38	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
acenaphthylene	< 0.38	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
fluorene	<b>0.43</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
anthracene	<b>0.84</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
fluoranthene	<b>6.8</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
pyrene	<b>5.8</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
benzo(a)anthracene	<b>3.0</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
chrysene	<b>3.6</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
benzo(b)fluoranthene	<b>2.8</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
benzo(k)fluoranthene	<b>2.7</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
benzo(a)pyrene	<b>3.0</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
indeno(1,2,3-cd)pyrene	<b>1.7</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
dibenzo(a,h)anthracene	<b>0.62</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
benzo(g,h,i)perylene	<b>2.0</b>	0.38	ug/g	1	CL	4/8/20	12630	4/9/20	19:14	MA EPH
Unadjusted C11-C22 Aromatics	<b>180</b>	38	ug/g	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
C9-C18 Aliphatics	< 38	38	ug/g	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
C19-C36 Aliphatics	<b>190</b>	38	ug/g	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
C11-C22 Aromatics	<b>140</b>	38	ug/g	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>52</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
o-terphenyl SUR	<b>59</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
2-fluorobiphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH
2-bromonaphthalene SUR	<b>64</b>	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:21	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-018

Sample ID: DUP-1

Matrix: Solid

Percent Dry: 23.7% Results expressed on a dry weight basis.

Sampled: 4/7/20 0:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	0.86	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
2-methylnaphthalene	< 0.83	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
phenanthrene	5.7	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
acenaphthene	< 0.83	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
acenaphthylene	< 0.83	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
fluorene	< 0.83	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
anthracene	0.92	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
fluoranthene	9.2	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
pyrene	9.0	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
benzo(a)anthracene	4.0	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
chrysene	5.3	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
benzo(b)fluoranthene	4.4	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
benzo(k)fluoranthene	3.7	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
benzo(a)pyrene	4.4	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
indeno(1,2,3-cd)pyrene	2.9	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
dibenzo(a,h)anthracene	< 0.83	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
benzo(g,h,i)perylene	3.2	0.83	ug/g	1	CL	4/8/20	12630	4/9/20	17:43	MA EPH
Unadjusted C11-C22 Aromatics	360	83	ug/g	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
C9-C18 Aliphatics	< 83	83	ug/g	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
C19-C36 Aliphatics	390	83	ug/g	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
C11-C22 Aromatics	300	83	ug/g	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	48	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
o-terphenyl SUR	54	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
2-fluorobiphenyl SUR	68	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH
2-bromonaphthalene SUR	66	40-140	%	1	DBV	4/8/20	12630	4/9/20	17:56	MA EPH

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-001

Sample ID: SB-104 (0-3')

Matrix: Solid Percent Dry: 89.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 8:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.54	0.54	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Arsenic	< 2.7	2.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Barium	17	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Beryllium	< 0.54	0.54	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Cadmium	< 0.54	0.54	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Chromium	< 5.4	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Lead	16	2.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/9/20	12632	4/15/20	12:18	SW7471B
Nickel	5.4	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Selenium	< 5.4	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Silver	< 2.7	2.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Thallium	< 0.54	0.54	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Vanadium	8.3	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A
Zinc	16	5.4	ug/g	5	AGN	4/9/20	12629	4/10/20	4:27	SW3051A6020A

Sample#: 52513-002

Sample ID: SB-104 (8-10')

Matrix: Solid Percent Dry: 83.1% Results expressed on a dry weight basis.

Sampled: 4/7/20 8:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.57	0.57	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Arsenic	3.1	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Barium	27	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Beryllium	< 0.57	0.57	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Cadmium	< 0.57	0.57	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Chromium	16	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Lead	5.5	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Mercury	< 0.16	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:20	SW7471B
Nickel	17	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Selenium	< 5.7	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Silver	< 2.8	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Thallium	< 0.57	0.57	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Vanadium	22	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A
Zinc	18	5.7	ug/g	5	AGN	4/9/20	12629	4/10/20	4:36	SW3051A6020A

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-004

Sample ID: SB-107 (0-3')

Matrix: Solid Percent Dry: 79% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	55	0.63	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Arsenic	8.0	3.2	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Barium	280 M	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
M = The recovery for the matrix spike was 156%. The acceptance criteria is 75-125%.										
Beryllium	< 0.63	0.63	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Cadmium	4.4	0.63	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Chromium	33	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Lead	1200 M	3.2	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
M = The recovery for the matrix spike was -7%. The acceptance criteria is 75-125%.										
Mercury	0.60 M	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:21	SW7471B
M = The recovery for the matrix spike/duplicate were 12% and 64%. The acceptance criteria is 80-120%.										
Nickel	76	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Selenium	< 6.3 M	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
M = The recovery for the matrix spike was 74%. The acceptance criteria is 75-125%.										
Silver	32	3.2	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Thallium	< 0.63	0.63	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Vanadium	20	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A
Zinc	2900	6.3	ug/g	5	AGN	4/9/20	12629	4/10/20	4:45	SW3051A6020A

Sample#: 52513-005

Sample ID: SB-107 (10-11')

Matrix: Solid Percent Dry: 84.5% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	5.9	0.49	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Arsenic	3.5	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Barium	48	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Beryllium	< 0.49	0.49	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Cadmium	2.0	0.49	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Chromium	19	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Lead	300	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Mercury	< 0.14	0.14	ug/g	1	AGN	4/9/20	12632	4/15/20	12:27	SW7471B
Nickel	16	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Selenium	< 4.9	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Silver	< 2.5	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Thallium	< 0.49	0.49	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Vanadium	12	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A
Zinc	250	4.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:13	SW3051A6020A

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-007

Sample ID: SB-108 (0-3')

Matrix: Solid Percent Dry: 73.6% Results expressed on a dry weight basis.

Sampled: 4/7/20 10:50

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	51	0.65	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Arsenic	8.1	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Barium	430	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Beryllium	< 0.65	0.65	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Cadmium	5.8	0.65	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Chromium	70	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Lead	1200	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Mercury	0.18	0.15	ug/g	1	AGN	4/9/20	12632	4/15/20	12:29	SW7471B
Nickel	88	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Selenium	< 6.5	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Silver	7.3	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Thallium	< 0.65	0.65	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Vanadium	10	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A
Zinc	2900	6.5	ug/g	5	AGN	4/9/20	12629	4/10/20	5:23	SW3051A6020A

Sample#: 52513-008

Sample ID: SB-108 (6-9')

Matrix: Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	10	0.60	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Arsenic	7.1	3.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Barium	470	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Beryllium	< 0.60	0.60	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Cadmium	11	0.60	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Chromium	25	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Lead	970	3.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Mercury	0.47	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:30	SW7471B
Nickel	40	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Selenium	< 6.0	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Silver	< 3.0	3.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Thallium	< 0.60	0.60	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Vanadium	18	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A
Zinc	2400	6.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:32	SW3051A6020A

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-010

Sample ID: SB-109 (0-3')

Matrix: Solid Percent Dry: 70.9% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	64	0.67	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Arsenic	23	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Barium	300	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Beryllium	< 0.67	0.67	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Cadmium	3.6	0.67	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Chromium	64	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Lead	1300	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Mercury	< 0.19	0.19	ug/g	1	AGN	4/9/20	12632	4/15/20	12:32	SW7471B
Nickel	100	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Selenium	< 6.7	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Silver	12	3.3	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Thallium	< 0.67	0.67	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Vanadium	8.1	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A
Zinc	4000	6.7	ug/g	5	AGN	4/9/20	12629	4/10/20	5:41	SW3051A6020A

Sample#: 52513-011

Sample ID: SB-109 (5-8')

Matrix: Solid Percent Dry: 23.7% Results expressed on a dry weight basis.

Sampled: 4/7/20 11:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	3.5	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Arsenic	15	9.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Barium	78	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Beryllium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Cadmium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Chromium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Lead	180	9.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Mercury	< 0.58	0.58	ug/g	1	AGN	4/9/20	12632	4/15/20	12:34	SW7471B
Nickel	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Selenium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Silver	< 9.9	9.9	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Thallium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Vanadium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A
Zinc	270	20	ug/g	5	AGN	4/9/20	12629	4/10/20	5:51	SW3051A6020A

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-013

Sample ID: SB-110 (0-3')

Matrix: Solid Percent Dry: 88.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 12:50

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.51	0.51	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Arsenic	<b>2.9</b>	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Barium	<b>41</b>	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Beryllium	< 0.51	0.51	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Cadmium	< 0.51	0.51	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Chromium	<b>10</b>	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Lead	<b>8.1</b>	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	19:48	SW3051A6020A
Mercury	< 0.16	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:36	SW7471B
Nickel	<b>12</b>	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Selenium	< 5.1	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Silver	< 2.5	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Thallium	< 0.51	0.51	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Vanadium	<b>16</b>	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A
Zinc	<b>19</b>	5.1	ug/g	5	AGN	4/9/20	12629	4/10/20	6:28	SW3051A6020A

Sample#: 52513-014

Sample ID: SB-110 (5-8')

Matrix: Solid Percent Dry: 17.7% Results expressed on a dry weight basis.

Sampled: 4/7/20 12:55

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	<b>3.5</b>	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Arsenic	< 14	14	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Barium	<b>160</b>	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Beryllium	< 2.8	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Cadmium	< 2.8	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Chromium	< 28	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Lead	<b>290</b>	14	ug/g	5	AGN	4/9/20	12629	4/10/20	19:56	SW3051A6020A
Mercury	< 0.78	0.78	ug/g	1	AGN	4/9/20	12632	4/15/20	12:41	SW7471B
Nickel	< 28	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Selenium	< 28	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Silver	< 14	14	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Thallium	< 2.8	2.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Vanadium	< 28	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A
Zinc	<b>230</b>	28	ug/g	5	AGN	4/9/20	12629	4/10/20	6:38	SW3051A6020A

Project ID: Cedar Woods 21

Job ID: 52513

Sample#: 52513-016

Sample ID: SB-112 (0-3')

Matrix: Solid Percent Dry: 90.3% Results expressed on a dry weight basis.

Sampled: 4/7/20 13:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.50	0.50	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Arsenic	<b>5.0</b>	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Barium	<b>38</b>	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Beryllium	< 0.50	0.50	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Cadmium	< 0.50	0.50	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Chromium	<b>18</b>	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Lead	<b>66</b>	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	20:04	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/9/20	12632	4/15/20	12:43	SW7471B
Nickel	<b>18</b>	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Selenium	< 5.0	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Silver	< 2.5	2.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Thallium	< 0.50	0.50	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Vanadium	<b>21</b>	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A
Zinc	<b>51</b>	5.0	ug/g	5	AGN	4/9/20	12629	4/10/20	6:47	SW3051A6020A

Sample#: 52513-017

Sample ID: SB-112 (5-8')

Matrix: Solid Percent Dry: 52.4% Results expressed on a dry weight basis.

Sampled: 4/7/20 13:25

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	<b>4.2</b>	0.95	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Arsenic	<b>16</b>	4.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Barium	<b>130</b>	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Beryllium	< 0.95	0.95	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Cadmium	<b>1.1</b>	0.95	ug/g	5	AGN	4/9/20	12629	4/10/20	20:13	SW3051A6020A
Chromium	<b>21</b>	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Lead	<b>470</b>	4.8	ug/g	5	AGN	4/9/20	12629	4/10/20	20:13	SW3051A6020A
Mercury	<b>0.80</b>	0.25	ug/g	1	AGN	4/9/20	12632	4/15/20	12:45	SW7471B
Nickel	<b>17</b>	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Selenium	< 9.5	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Silver	< 4.8	4.8	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Thallium	< 0.95	0.95	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Vanadium	<b>26</b>	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A
Zinc	<b>300</b>	9.5	ug/g	5	AGN	4/9/20	12629	4/10/20	6:56	SW3051A6020A



**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-018

**Sample ID:** DUP-1

**Matrix:** Solid

Percent Dry: 23.7% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 0:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	<b>2.1</b>	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Arsenic	<b>20</b>	10	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Barium	<b>86</b>	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Beryllium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Cadmium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Chromium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Lead	<b>250</b>	10	ug/g	5	AGN	4/9/20	12629	4/10/20	20:46	SW3051A6020A
Mercury	<b>0.99</b>	0.54	ug/g	1	AGN	4/9/20	12632	4/15/20	12:47	SW7471B
Nickel	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Selenium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Silver	< 10	10	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Thallium	< 2.0	2.0	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Vanadium	< 20	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A
Zinc	<b>250</b>	20	ug/g	5	AGN	4/9/20	12629	4/10/20	7:06	SW3051A6020A

**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-002

**Sample ID:** SB-104 (8-10')

**Matrix:** Solid Percent Dry: 83.1% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 8:55

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.47	0.47	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	<b>470</b>		mV	1	WAS	2001839	4/8/20	6:07	ASTM1498-08	
pH	<b>5.9 D</b>		pH	1	WAS	2001840	4/8/20	5:55	SW9045C	

D = The difference for the sample duplicate, run as internal QC, was outside the 0.2 pH units acceptance range. The duplicate result is 5.6 pH units.

**Sample#:** 52513-004

**Sample ID:** SB-107 (0-3')

**Matrix:** Solid Percent Dry: 79% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 10:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.51 M	0.51	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	<b>180</b>		mV	1	WAS	2001839	4/8/20	6:28	ASTM1498-08	
pH	<b>8.3</b>		pH	1	WAS	2001840	4/8/20	6:00	SW9045C	

M = The percent recovery in the matrix spike was outside acceptance criteria. See case narrative.

**Sample#:** 52513-008

**Sample ID:** SB-108 (6-9')

**Matrix:** Solid Percent Dry: 80.4% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 11:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.48	0.48	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	<b>200</b>		mV	1	WAS	2001839	4/8/20	6:35	ASTM1498-08	
pH	<b>8.0</b>		pH	1	WAS	2001840	4/8/20	6:08	SW9045C	

**Sample#:** 52513-011

**Sample ID:** SB-109 (5-8')

**Matrix:** Solid Percent Dry: 23.7% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 11:45

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 1.6	1.6	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	<b>140</b>		mV	1	WAS	2001839	4/8/20	6:48	ASTM1498-08	
pH	<b>7.5</b>		pH	1	WAS	2001840	4/8/20	6:18	SW9045C	

**Sample#:** 52513-014

**Sample ID:** SB-110 (5-8')

**Matrix:** Solid Percent Dry: 17.7% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 12:55

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 2.2	2.2	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	<b>130</b>		mV	1	WAS	2001839	4/8/20	7:02	ASTM1498-08	
pH	<b>7.4</b>		pH	1	WAS	2001840	4/8/20	6:20	SW9045C	

**Project ID:** Cedar Woods 21

**Job ID:** 52513

**Sample#:** 52513-017

**Sample ID:** SB-112 (5-8')

**Matrix:** Solid      Percent Dry: 52.4% Results expressed on a dry weight basis.

**Sampled:** 4/7/20    13:25

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Chromium, Hexavalent	1.3	0.76	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	140		mV	1	WAS	2001839	4/8/20	7:10	ASTM1498-08	
pH	7.6		pH	1	WAS	2001840	4/8/20	6:22	SW9045C	

# Quality Control Report



124 Heritage Avenue Unit 16  
Portsmouth, NH 03801

[www.absoluteresourceassociates.com](http://www.absoluteresourceassociates.com)

## MassDEP Analytical Protocol Certification Form

Laboratory Name: Absolute Resource Associates

Project #: 21

Project Location: Massachusetts

RTN:

**This Form provides certifications for the following data set: list Laboratory Sample ID Number(s): 52513**

Matrices:  Groundwater/Surface Water  Soil/Sediment  Drinking Water  Air  Other:

**CAM Protocol** (check all that apply below):

8260 VOC CAM II A <input checked="" type="checkbox"/>	7470/7471 Hg CAM III B <input checked="" type="checkbox"/>	MassDEP VPH (GC/PID/FID) CAM IV A <input checked="" type="checkbox"/>	8082 PCB CAM V A <input type="checkbox"/>	9014 Total Cyanide/PAC CAM VI A <input type="checkbox"/>	6860 Perchlorate CAM VIII B <input type="checkbox"/>
8270 SVOC CAM II B <input type="checkbox"/>	7010 Metals CAM III C <input type="checkbox"/>	MassDEP VPH (GC/MS) CAM IV C <input type="checkbox"/>	8081 Pesticides CAM V B <input type="checkbox"/>	7196 Hex Cr CAM VI B <input checked="" type="checkbox"/>	MassDEP APH CAM IX A <input type="checkbox"/>
6010 Metals CAM III A <input type="checkbox"/>	6020 Metals CAM III D <input checked="" type="checkbox"/>	MassDEP EPH CAM IV B <input checked="" type="checkbox"/>	8151 Herbicides CAM V C <input type="checkbox"/>	8330 Explosives CAM VIII A <input type="checkbox"/>	TO-15 VOC CAM IX B <input type="checkbox"/>

**Affirmative Responses to Questions A through F are required for "Presumptive Certainty" status**

<b>A</b>	Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature) in the field or laboratory, and prepared/analyzed within method holding times?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>B</b>	Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>C</b>	Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>D</b>	Does the laboratory report comply with all the reporting requirements specified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>E</b>	VPH, EPH, APH, and TO-15 only a. VPH, EPH, and APH Methods only: Was each method conducted without significant modification(s)? (Refer to the individual method(s) for a list of significant modifications). b. APH and TO-15 Methods only: Was the complete analyte list reported for each method?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Yes <input type="checkbox"/> No
<b>F</b>	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No

**Responses to Questions G, H and I below are required for "Presumptive Certainty" status**

<b>G</b>	Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <sup>1</sup>
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**Data User Note:** Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40. 1056 (2)(k) and WSC-07-350.

<b>H</b>	Were all QC performance standards specified in the CAM protocol(s) achieved?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
<b>I</b>	Were results reported for the complete analyte list specified in the selected CAM protocol(s)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <sup>1</sup>

<sup>1</sup>All negative responses must be addressed in an attached laboratory narrative.

*I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, is accurate and complete.*

Signature: 

Position: Chief Operating Officer

Printed Name: Aaron DeWees

Date: 5/1/20

## Sample Integrity Table

Parameter	Method	Matrix	Minimum Volume	Recommended Container(s)	Required Preservation	Holding Time
Volatile Organics	EPA 8260	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
Volatile Organics	EPA 8260	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	14 Days
Semivolatile Organics	EPA 8270	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Semivolatile Organics	EPA 8270	Solid	20g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Organochlorine Pesticides	EPA 8081	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Organochlorine Pesticides	EPA 8081	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
PCBs	EPA 8082	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
PCBs	EPA 8082	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
Herbicides (subcontracted)	EPA 8151	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Herbicides (subcontracted)	EPA 8151	Solid	30g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
MA DEP VPH	MADEP VPH	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP VPH	MADEP VPH	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	28 Days
MA DEP EPH	MADEP EPH	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP EPH	MADEP EPH	Solid	30g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Total Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	180 Days
Dissolved Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	Filter First 1:1 HNO <sub>3</sub> to pH <2	180 Days
Total Metals	EPA 6010	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	180 Days
Total Mercury (may be combined with Total Metals)	EPA 7470	Aqueous	100mL	125mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	28 Days
Total Mercury (may be combined with Total Metals)	EPA 7471	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	28 Days
Chromium, Hexavalent	EPA 7196	Aqueous	100mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> buffer	28 Days
Chromium, Hexavalent (subcontract)	EPA 7196	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	30 Days
Cyanide, Total	EPA 9014	Aqueous	125mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ 1:1 NaOH to pH >8	14 Days
Cyanide, Total	EPA 9014	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days

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## Case Narrative

Lab # 52513

### **Sample Receiving and Chain of Custody Discrepancies**

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Samples were received in acceptable condition, between 0 and 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

As noted on the result page, several VOC samples did not meet the 1:1 +/-25% methanol to soil ratio.

### **Calibration**

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VOC: See the included table for a list of compounds quantitated by quadratic equation.

SVOC: See the included table for a list of compounds quantitated by quadratic equation.

### **Method Blank**

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No exceptions noted.

### **Surrogate Recoveries**

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VPH: Sample 52513-004 did not meet acceptance criteria for the surrogate a,a,a-trifluorotoluene. Matrix interference suspected. Sample 52513-005 did not meet acceptance criteria for the aliphatic surrogate 2,5-dibromotoluene, likely as a result of hydrocarbon interference. Sample 52513-014 did not meet acceptance criteria for the surrogate a,a,a-trifluorotoluene, likely as a result of the sample's high moisture content. All of the associated sample chromatograms are included in the report.

VOC: The LCS/D12627 did not meet acceptance limits for the surrogate 4-bromofluorobenzene. The percent recovery for this surrogate was outside of the acceptance criteria in sample 52513-008, likely as a result of hydrocarbon interference. The sample chromatogram is included in the report.

### **Laboratory Control Sample Results**

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VOC: The LCS12627 did not meet the acceptance criteria for bromomethane and t-butanol (TBA). Since recoveries were >10% and <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

### **Matrix Spike/Matrix Spike Duplicate/Duplicate Results**

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Metals: The percent recovery for mercury in the MS/D12632 (52513-004) was 12% and 64% respectively, outside the acceptance criteria of 80-120%. All other batch QC was within acceptance limits. Matrix interference is suspected. The percent recovery for lead in the MS/D12629 was -7% and 43% respectively, outside the acceptance criteria of 75-125%. All other batch QC was within acceptance limits. Matrix interference is suspected. The percent recovery in the MSD12629 was 156% for barium and 74% for selenium. All other batch QC was within acceptance limits. Matrix interference is suspected. All samples with results greater than the reporting limit for the affected compounds have been qualified accordingly.



**Case Narrative**

**Lab # 52513**

Chromium, Hexavalent: The MS/D2001871 (52513-004) did not meet the acceptance criteria. The percent recovery was acceptable in the associated CRM/CRMD. The results for the ORP and pH analysis show that the sample exhibits a reducing nature therefore resulting in poor spike recovery. The sample has been qualified accordingly.

**Other**

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No other exceptions noted.

VPH: The trap used for VPH analysis is a Tekmar STRATUM Purge Trap 9. The column used for VPH analysis is a Restek Rtx-502.2, 105m, 0.53mmID, and 3um df.

**MassDEP Analytical Protocol Certification Form Questions A through I**

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No explanation is needed for Questions A through I answered in the affirmative.

**Question H:** See surrogate section above. Box H is "No."





**Quantitation by Quadratic Equation**  
**Lab # 52513**

VOC: Quantitation of the following compounds was based on a quadratic equation:

Acetone

2-Hexanone

Bromoform

hexachlorobutadiene

SVOC: Quantitation of the following compounds was based on a quadratic equation:

2-methylnaphthalene

benzo(k)fluoranthene

indeno(1,2,3-cd)pyrene

dibenzo(a,h)anthracene

benzo(g,h,i)perylene

# GLOSSARY

%R	Percent Recovery
BLK	Blank (Method Blank, Preparation Blank)
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRM	Certified Reference Material (associated with solid Metals samples)
CRMD	Certified Reference Material Duplicate (associated with solid Metals samples)
Dil'n	Dilution
DL	Detection Limit
DUP	Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
LOD	Limit of Detection
LOQ	Limit of Quantitation
MB	Methanol Blank (associated with solid VOC samples)
MLCS	Methanol Laboratory Control Sample (associated with solid VOC samples)
MLCSD	Methanol Laboratory Control Sample Duplicate (associated with solid VOC samples)
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PB	Preparation Blank
QC	Quality Control
RL	Reporting Limit
RPD	Relative Percent Difference
SUR	Surrogate



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- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
MA VPH	MB12628	Unadjusted C5-C8 Aliphatics	<	5.0	ug/g						
		Unadjusted C9-C12 Aliphatics	<	5.0	ug/g						
		methyl t-butyl ether (MTBE)	<	0.10	ug/g						
		benzene	<	0.10	ug/g						
		toluene	<	0.10	ug/g						
		ethylbenzene	<	0.10	ug/g						
		m&p-xylenes	<	0.10	ug/g						
		o-xylene	<	0.10	ug/g						
		naphthalene	<	0.25	ug/g						
		C5-C8 Aliphatics	<	5.0	ug/g						
		C9-C12 Aliphatics	<	5.0	ug/g						
		C9-C10 Aromatics	<	5.0	ug/g						
		2,5-dibromotoluene as Aromatic SUR		113	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		
MA VPH	MLCS12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130		
		Unadjusted C9-C12 Aliphatics		13	ug/g	15	89	70	130		
		methyl t-butyl ether (MTBE)		4.6	ug/g	5	92	70	130		
		benzene		4.8	ug/g	5	97	70	130		
		toluene		4.8	ug/g	5	95	70	130		
		ethylbenzene		4.8	ug/g	5	96	70	130		
		m&p-xylenes		9.7	ug/g	10	97	70	130		
		o-xylene		4.8	ug/g	5	96	70	130		
		naphthalene		5.0	ug/g	5	100	70	130		
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		
		C9-C10 Aromatics		5.1	ug/g	5	102	70	130		
		2,5-dibromotoluene as Aromatic SUR		111	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		86	%			70	130		
MA VPH	MLCSD12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130	0	25
		Unadjusted C9-C12 Aliphatics		14	ug/g	15	90	70	130	2	25
		methyl t-butyl ether (MTBE)		4.7	ug/g	5	93	70	130	1	25
		benzene		4.9	ug/g	5	99	70	130	2	25
		toluene		4.9	ug/g	5	98	70	130	3	25
		ethylbenzene		4.9	ug/g	5	99	70	130	3	25
		m&p-xylenes		10.0	ug/g	10	100	70	130	3	25
		o-xylene		4.9	ug/g	5	99	70	130	2	25
		naphthalene		5.1	ug/g	5	102	70	130	2	25
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C10 Aromatics		5.2	ug/g	5	103	70	130	1	25
		2,5-dibromotoluene as Aromatic SUR		99	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		100	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MB12627	dichlorodifluoromethane		<	0.10	ug/g				
		chloromethane		<	0.10	ug/g				
		vinyl chloride		<	0.10	ug/g				
		bromomethane		<	0.25	ug/g				
		chloroethane		<	0.10	ug/g				
		trichlorofluoromethane		<	0.10	ug/g				
		diethyl ether		<	0.50	ug/g				
		acetone		<	2.5	ug/g				
		1,1-dichloroethene		<	0.10	ug/g				
		methylene chloride		<	0.25	ug/g				
		carbon disulfide		<	0.10	ug/g				
		methyl t-butyl ether (MTBE)		<	0.10	ug/g				
		trans-1,2-dichloroethene		<	0.10	ug/g				
		isopropyl ether (DIPE)		<	0.10	ug/g				
		ethyl t-butyl ether (ETBE)		<	0.10	ug/g				
		1,1-dichloroethane		<	0.10	ug/g				
		t-butanol (TBA)		<	2.5	ug/g				
		2-butanone (MEK)		<	0.30	ug/g				
		2,2-dichloropropane		<	0.10	ug/g				
		cis-1,2-dichloroethene		<	0.10	ug/g				
		chloroform		<	0.10	ug/g				
		bromochloromethane		<	0.10	ug/g				
		tetrahydrofuran (THF)		<	0.50	ug/g				
		1,1,1-trichloroethane		<	0.10	ug/g				
		1,1-dichloropropene		<	0.10	ug/g				
		t-amyl-methyl ether (TAME)		<	0.10	ug/g				
		carbon tetrachloride		<	0.10	ug/g				
		1,2-dichloroethane		<	0.10	ug/g				
		benzene		<	0.10	ug/g				
		trichloroethene		<	0.10	ug/g				
		1,2-dichloropropane		<	0.10	ug/g				
		bromodichloromethane		<	0.10	ug/g				
		1,4-dioxane		<	2.5	ug/g				
		dibromomethane		<	0.10	ug/g				
		4-methyl-2-pentanone (MIBK)		<	0.45	ug/g				
		cis-1,3-dichloropropene		<	0.10	ug/g				
		toluene		<	0.10	ug/g				
		trans-1,3-dichloropropene		<	0.10	ug/g				
		2-hexanone		<	0.50	ug/g				
		1,1,2-trichloroethane		<	0.10	ug/g				
		1,3-dichloropropane		<	0.10	ug/g				
		tetrachloroethene		<	0.10	ug/g				
		dibromochloromethane		<	0.10	ug/g				
		1,2-dibromoethane (EDB)		<	0.10	ug/g				
		chlorobenzene		<	0.10	ug/g				
		1,1,1,2-tetrachloroethane		<	0.10	ug/g				
		ethylbenzene		<	0.10	ug/g				
		m&p-xylenes		<	0.10	ug/g				
		o-xylene		<	0.10	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MB12627	styrene		<	0.10	ug/g				
		bromoform		<	0.10	ug/g				
		isopropylbenzene		<	0.10	ug/g				
		1,1,2,2-tetrachloroethane		<	0.10	ug/g				
		1,2,3-trichloropropane		<	0.10	ug/g				
		n-propylbenzene		<	0.10	ug/g				
		bromobenzene		<	0.10	ug/g				
		1,3,5-trimethylbenzene		<	0.10	ug/g				
		2-chlorotoluene		<	0.10	ug/g				
		4-chlorotoluene		<	0.10	ug/g				
		tert-butylbenzene		<	0.10	ug/g				
		1,2,4-trimethylbenzene		<	0.10	ug/g				
		sec-butylbenzene		<	0.10	ug/g				
		1,3-dichlorobenzene		<	0.10	ug/g				
		4-isopropyltoluene		<	0.10	ug/g				
		1,4-dichlorobenzene		<	0.10	ug/g				
		1,2-dichlorobenzene		<	0.10	ug/g				
		n-butylbenzene		<	0.10	ug/g				
		1,2-dibromo-3-chloropropane (DBCP)		<	0.10	ug/g				
		1,2,4-trichlorobenzene		<	0.10	ug/g				
		hexachlorobutadiene		<	0.10	ug/g				
		naphthalene		<	0.25	ug/g				
		1,2,3-trichlorobenzene		<	0.10	ug/g				
		dibromofluoromethane SUR			100	%		78	114	
		toluene-D8 SUR			104	%		88	110	
		4-bromofluorobenzene SUR			105	%		86	115	
		a,a,a-trifluorotoluene SUR			122	%		70	130	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MLCS12627	dichlorodifluoromethane		0.79	ug/g	1	79	70	130	
		chloromethane		0.95	ug/g	1	95	70	130	
		vinyl chloride		0.93	ug/g	1	93	70	130	
		bromomethane		0.64	ug/g	1	64	70	130	*
		chloroethane		1.0	ug/g	1	104	70	130	
		trichlorofluoromethane		1.1	ug/g	1	106	70	130	
		diethyl ether		0.99	ug/g	1	99	70	130	
		acetone	<	2.5	ug/g	1	120			
		1,1-dichloroethene		0.97	ug/g	1	97	70	130	
		methylene chloride		1.1	ug/g	1	106	70	130	
		carbon disulfide		0.94	ug/g	1	94	70	130	
		methyl t-butyl ether (MTBE)		0.96	ug/g	1	96	70	130	
		trans-1,2-dichloroethene		1.0	ug/g	1	102	70	130	
		isopropyl ether (DIPE)		1.0	ug/g	1	101	70	130	
		ethyl t-butyl ether (ETBE)		0.98	ug/g	1	98	70	130	
		1,1-dichloroethane		1.0	ug/g	1	103	70	130	
		t-butanol (TBA)		6.6	ug/g	5	132	70	130	*
		2-butanone (MEK)		1.2	ug/g	1	119	70	130	
		2,2-dichloropropane		0.90	ug/g	1	90	70	130	
		cis-1,2-dichloroethene		1.1	ug/g	1	106	70	130	
		chloroform		1.0	ug/g	1	102	70	130	
		bromochloromethane		0.98	ug/g	1	98	70	130	
		tetrahydrofuran (THF)		1.1	ug/g	1	106	70	130	
		1,1,1-trichloroethane		0.99	ug/g	1	99	70	130	
		1,1-dichloropropene		0.95	ug/g	1	95	70	130	
		t-amyl-methyl ether (TAME)		0.92	ug/g	1	92	70	130	
		carbon tetrachloride		0.98	ug/g	1	98	70	130	
		1,2-dichloroethane		0.97	ug/g	1	97	70	130	
		benzene		0.99	ug/g	1	99	70	130	
		trichloroethene		1.0	ug/g	1	101	70	130	
		1,2-dichloropropane		0.99	ug/g	1	99	70	130	
		bromodichloromethane		1.00	ug/g	1	100	70	130	
		1,4-dioxane	<	2.5	ug/g	2	117	70	130	
		dibromomethane		1.1	ug/g	1	105	70	130	
		4-methyl-2-pentanone (MIBK)		1.1	ug/g	1	109	70	130	
		cis-1,3-dichloropropene		1.0	ug/g	1	101	70	130	
		toluene		0.99	ug/g	1	99	70	130	
		trans-1,3-dichloropropene		0.98	ug/g	1	98	70	130	
		2-hexanone		1.2	ug/g	1	117	70	130	
		1,1,2-trichloroethane		1.1	ug/g	1	107	70	130	
		1,3-dichloropropane		1.0	ug/g	1	102	70	130	
		tetrachloroethene		1.0	ug/g	1	103	70	130	
		dibromochloromethane		0.98	ug/g	1	98	70	130	
		1,2-dibromoethane (EDB)		1.0	ug/g	1	104	70	130	
		chlorobenzene		1.0	ug/g	1	103	70	130	
		1,1,1,2-tetrachloroethane		1.0	ug/g	1	102	70	130	
		ethylbenzene		1.0	ug/g	1	100	70	130	
		m&p-xylenes		2.1	ug/g	2	107	70	130	
		o-xylene		1.1	ug/g	1	106	70	130	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MLCS12627	styrene		1.1	ug/g	1	106	70	130	
		bromoform		1.1	ug/g	1	110	70	130	
		isopropylbenzene		1.1	ug/g	1	108	70	130	
		1,1,2,2-tetrachloroethane		1.0	ug/g	1	104	70	130	
		1,2,3-trichloropropane		1.1	ug/g	1	107	70	130	
		n-propylbenzene		0.98	ug/g	1	98	70	130	
		bromobenzene		1.1	ug/g	1	106	70	130	
		1,3,5-trimethylbenzene		1.1	ug/g	1	107	70	130	
		2-chlorotoluene		1.0	ug/g	1	102	70	130	
		4-chlorotoluene		1.1	ug/g	1	106	70	130	
		tert-butylbenzene		0.97	ug/g	1	97	70	130	
		1,2,4-trimethylbenzene		1.0	ug/g	1	101	70	130	
		sec-butylbenzene		0.95	ug/g	1	95	70	130	
		1,3-dichlorobenzene		1.1	ug/g	1	106	70	130	
		4-isopropyltoluene		1.0	ug/g	1	104	70	130	
		1,4-dichlorobenzene		1.1	ug/g	1	107	70	130	
		1,2-dichlorobenzene		1.1	ug/g	1	107	70	130	
		n-butylbenzene		1.0	ug/g	1	101	70	130	
		1,2-dibromo-3-chloropropane (DBCP)		1.1	ug/g	1	112	70	130	
		1,2,4-trichlorobenzene		1.0	ug/g	1	103	70	130	
		hexachlorobutadiene		1.1	ug/g	1	107	70	130	
		naphthalene		1.1	ug/g	1	109	70	130	
		1,2,3-trichlorobenzene		1.1	ug/g	1	108	70	130	
		dibromofluoromethane SUR		87	%			78	114	
		toluene-D8 SUR		104	%			88	110	
		4-bromofluorobenzene SUR		118	%		*	86	115	
		a,a,a-trifluorotoluene SUR		102	%			70	130	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MLCSD12627	dichlorodifluoromethane		0.82	ug/g	1	82	70 130	4	30
		chloromethane		1.0	ug/g	1	101	70 130	6	30
		vinyl chloride		0.94	ug/g	1	94	70 130	1	30
		bromomethane		0.86	ug/g	1	86	70 130	30	30
		chloroethane		1.1	ug/g	1	107	70 130	3	30
		trichlorofluoromethane		1.1	ug/g	1	108	70 130	3	30
		diethyl ether		1.0	ug/g	1	100	70 130	1	30
		acetone	<	2.5	ug/g	1	109		9	30
		1,1-dichloroethene		0.98	ug/g	1	98	70 130	1	30
		methylene chloride		1.1	ug/g	1	111	70 130	4	30
		carbon disulfide		0.96	ug/g	1	96	70 130	1	30
		methyl t-butyl ether (MTBE)		0.97	ug/g	1	97	70 130	1	30
		trans-1,2-dichloroethene		1.0	ug/g	1	104	70 130	3	30
		isopropyl ether (DIPE)		1.0	ug/g	1	101	70 130	0	30
		ethyl t-butyl ether (ETBE)		1.0	ug/g	1	100	70 130	2	30
		1,1-dichloroethane		1.0	ug/g	1	103	70 130	0	30
		t-butanol (TBA)		6.1	ug/g	5	122	70 130	8	30
		2-butanone (MEK)		1.1	ug/g	1	113	70 130	5	30
		2,2-dichloropropane		0.91	ug/g	1	91	70 130	2	30
		cis-1,2-dichloroethene		1.1	ug/g	1	107	70 130	1	30
		chloroform		1.1	ug/g	1	106	70 130	3	30
		bromochloromethane		1.0	ug/g	1	103	70 130	5	30
		tetrahydrofuran (THF)		1.0	ug/g	1	102	70 130	4	30
		1,1,1-trichloroethane		1.0	ug/g	1	100	70 130	1	30
		1,1-dichloropropene		1.00	ug/g	1	100	70 130	4	30
		t-amyl-methyl ether (TAME)		0.93	ug/g	1	93	70 130	1	30
		carbon tetrachloride		0.99	ug/g	1	99	70 130	2	30
		1,2-dichloroethane		1.00	ug/g	1	100	70 130	3	30
		benzene		1.0	ug/g	1	101	70 130	3	30
		trichloroethene		1.0	ug/g	1	103	70 130	2	30
		1,2-dichloropropane		1.0	ug/g	1	101	70 130	1	30
		bromodichloromethane		1.0	ug/g	1	101	70 130	2	30
		1,4-dioxane	<	2.5	ug/g	2	115	70 130	2	30
		dibromomethane		1.1	ug/g	1	106	70 130	1	30
		4-methyl-2-pentanone (MIBK)		1.1	ug/g	1	105	70 130	3	30
		cis-1,3-dichloropropene		1.0	ug/g	1	103	70 130	2	30
		toluene		1.0	ug/g	1	101	70 130	2	30
		trans-1,3-dichloropropene		1.0	ug/g	1	101	70 130	3	30
		2-hexanone		1.1	ug/g	1	109	70 130	7	30
		1,1,2-trichloroethane		1.1	ug/g	1	110	70 130	3	30
		1,3-dichloropropane		1.0	ug/g	1	101	70 130	0	30
		tetrachloroethene		1.0	ug/g	1	103	70 130	0	30
		dibromochloromethane		1.00	ug/g	1	100	70 130	2	30
		1,2-dibromoethane (EDB)		1.0	ug/g	1	103	70 130	1	30
		chlorobenzene		1.0	ug/g	1	103	70 130	1	30
		1,1,1,2-tetrachloroethane		1.0	ug/g	1	102	70 130	1	30
		ethylbenzene		1.00	ug/g	1	100	70 130	1	30
		m&p-xylenes		2.1	ug/g	2	107	70 130	1	30
		o-xylene		1.1	ug/g	1	107	70 130	1	30



Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260D	MLCSD12627	styrene		1.1	ug/g	1	107	70 130	1	30
		bromoform		1.1	ug/g	1	111	70 130	1	30
		isopropylbenzene		1.1	ug/g	1	109	70 130	1	30
		1,1,2,2-tetrachloroethane		1.0	ug/g	1	102	70 130	2	30
		1,2,3-trichloropropane		1.0	ug/g	1	104	70 130	2	30
		n-propylbenzene		0.98	ug/g	1	98	70 130	0	30
		bromobenzene		1.0	ug/g	1	104	70 130	2	30
		1,3,5-trimethylbenzene		1.1	ug/g	1	106	70 130	1	30
		2-chlorotoluene		1.0	ug/g	1	101	70 130	1	30
		4-chlorotoluene		1.1	ug/g	1	107	70 130	1	30
		tert-butylbenzene		0.95	ug/g	1	95	70 130	2	30
		1,2,4-trimethylbenzene		1.0	ug/g	1	101	70 130	0	30
		sec-butylbenzene		0.94	ug/g	1	94	70 130	1	30
		1,3-dichlorobenzene		1.1	ug/g	1	106	70 130	0	30
		4-isopropyltoluene		1.0	ug/g	1	104	70 130	0	30
		1,4-dichlorobenzene		1.1	ug/g	1	107	70 130	0	30
		1,2-dichlorobenzene		1.1	ug/g	1	106	70 130	1	30
		n-butylbenzene		1.00	ug/g	1	100	70 130	2	30
		1,2-dibromo-3-chloropropane (DBCP)		1.0	ug/g	1	105	70 130	6	30
		1,2,4-trichlorobenzene		1.0	ug/g	1	102	70 130	1	30
		hexachlorobutadiene		1.0	ug/g	1	104	70 130	2	30
		naphthalene		1.0	ug/g	1	103	70 130	6	30
		1,2,3-trichlorobenzene		1.0	ug/g	1	103	70 130	4	30
		dibromofluoromethane SUR		88	%			78 114		
		toluene-D8 SUR		106	%			88 110		
		4-bromofluorobenzene SUR		116	%		*	86 115		
		a,a,a-trifluorotoluene SUR		109	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK12630	naphthalene		<	0.20	ug/g				
		2-methylnaphthalene		<	0.20	ug/g				
		phenanthrene		<	0.20	ug/g				
		acenaphthene		<	0.20	ug/g				
		acenaphthylene		<	0.20	ug/g				
		fluorene		<	0.20	ug/g				
		anthracene		<	0.20	ug/g				
		fluoranthene		<	0.20	ug/g				
		pyrene		<	0.20	ug/g				
		benzo(a)anthracene		<	0.20	ug/g				
		chrysene		<	0.20	ug/g				
		benzo(b)fluoranthene		<	0.20	ug/g				
		benzo(k)fluoranthene		<	0.20	ug/g				
		benzo(a)pyrene		<	0.20	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.20	ug/g				
		dibenzo(a,h)anthracene		<	0.20	ug/g				
		benzo(g,h,i)perylene		<	0.20	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			52	%		40	140	
		o-terphenyl SUR			59	%		40	140	
		2-fluorobiphenyl SUR			66	%		40	140	
		2-bromonaphthalene SUR			64	%		40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCS12630	naphthalene		3.1	ug/g	6	51	40	140	
		2-methylnaphthalene		3.1	ug/g	6	52	40	140	
		phenanthrene		3.6	ug/g	6	60	40	140	
		acenaphthene		3.3	ug/g	6	56	40	140	
		acenaphthylene		3.3	ug/g	6	54	40	140	
		fluorene		3.6	ug/g	6	60	40	140	
		anthracene		3.5	ug/g	6	59	40	140	
		fluoranthene		4.0	ug/g	6	66	40	140	
		pyrene		4.2	ug/g	6	70	40	140	
		benzo(a)anthracene		4.3	ug/g	6	71	40	140	
		chrysene		4.4	ug/g	6	73	40	140	
		benzo(b)fluoranthene		4.4	ug/g	6	73	40	140	
		benzo(k)fluoranthene		3.9	ug/g	6	65	40	140	
		benzo(a)pyrene		4.1	ug/g	6	68	40	140	
		indeno(1,2,3-cd)pyrene		4.4	ug/g	6	73	40	140	
		dibenzo(a,h)anthracene		4.4	ug/g	6	73	40	140	
		benzo(g,h,i)perylene		4.4	ug/g	6	73	40	140	
		Unadjusted C11-C22 Aromatics		65	ug/g	102	64	40	140	
		C9-C18 Aliphatics	<	20	ug/g	36	44	40	140	
		C19-C36 Aliphatics		49	ug/g	48	102	40	140	
		C11-C22 Aromatics	<	20	ug/g			40	140	
		1-chloro-octadecane SUR		52	%			40	140	
		o-terphenyl SUR		57	%			40	140	
		2-fluorobiphenyl SUR		64	%			40	140	
		2-bromonaphthalene SUR		61	%			40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD12630	naphthalene		2.9	ug/g	6	48	40 140	6	25
		2-methylnaphthalene		2.9	ug/g	6	49	40 140	6	25
		phenanthrene		3.5	ug/g	6	59	40 140	2	25
		acenaphthene		3.3	ug/g	6	55	40 140	1	25
		acenaphthylene		3.1	ug/g	6	52	40 140	4	25
		fluorene		3.6	ug/g	6	59	40 140	1	25
		anthracene		3.5	ug/g	6	59	40 140	0	25
		fluoranthene		4.1	ug/g	6	68	40 140	3	25
		pyrene		4.0	ug/g	6	66	40 140	5	25
		benzo(a)anthracene		4.2	ug/g	6	70	40 140	2	25
		chrysene		4.3	ug/g	6	71	40 140	2	25
		benzo(b)fluoranthene		4.1	ug/g	6	68	40 140	8	25
		benzo(k)fluoranthene		4.2	ug/g	6	70	40 140	8	25
		benzo(a)pyrene		4.1	ug/g	6	68	40 140	0	25
		indeno(1,2,3-cd)pyrene		4.2	ug/g	6	71	40 140	4	25
		dibenzo(a,h)anthracene		4.2	ug/g	6	71	40 140	3	25
		benzo(g,h,i)perylene		4.1	ug/g	6	69	40 140	6	25
		Unadjusted C11-C22 Aromatics		67	ug/g	102	66	40 140	3	25
		C9-C18 Aliphatics	<	20	ug/g	36	53	40 140	18	25
		C19-C36 Aliphatics		54	ug/g	48	112	40 140	9	25
		C11-C22 Aromatics	<	20	ug/g			40 140		25
		1-chloro-octadecane SUR		63	%			40 140		
		o-terphenyl SUR		60	%			40 140		
		2-fluorobiphenyl SUR		60	%			40 140		
		2-bromonaphthalene SUR		59	%			40 140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3051A6020A	BLK12629	Silver		< 0.025	mg/L						
		Arsenic		< 0.025	mg/L						
		Barium		< 0.050	mg/L						
		Beryllium		< 0.0050	mg/L						
		Cadmium		< 0.0050	mg/L						
		Chromium		< 0.050	mg/L						
		Nickel		< 0.050	mg/L						
		Lead		< 0.025	mg/L						
		Antimony		< 0.0050	mg/L						
		Selenium		< 0.050	mg/L						
		Thallium		< 0.0050	mg/L						
		Vanadium		< 0.050	mg/L						
		Zinc		< 0.050	mg/L						
SW3051A6020A	CRM12629	Silver		44.2	ug/g	53.8		31.9	59.7		
		Arsenic		153	ug/g	219		129	240		
		Barium		586	ug/g	788		509	867		
		Beryllium		183	ug/g	247		160	272		
		Cadmium		141	ug/g	175		111	192		
		Chromium		273	ug/g	375		223	414		
		Nickel		232	ug/g	318		193	358		
		Lead		254	ug/g	321		207	353		
		Antimony		82.9	ug/g	159		15.9	180		
		Selenium		97.8	ug/g	145		78.5	160		
		Thallium		80.5	ug/g	98.6		53.7	111		
		Vanadium		182	ug/g	267		168	294		
		Zinc		215	ug/g	311		190	352		
SW3051A6020A	CRMD12629	Silver		45.1	ug/g	53.8		31.9	59.7	2	20
		Arsenic		157	ug/g	219		129	240	3	20
		Barium		567	ug/g	788		509	867	3	20
		Beryllium		181	ug/g	247		160	272	1	20
		Cadmium		141	ug/g	175		111	192	0	20
		Chromium		278	ug/g	375		223	414	2	20
		Nickel		238	ug/g	318		193	358	3	20
		Lead		250	ug/g	321		207	353	2	20
		Antimony		82.8	ug/g	159		15.9	180	0	20
		Selenium		100	ug/g	145		78.5	160	2	20
		Thallium		80.8	ug/g	98.6		53.7	111	0	20
		Vanadium		185	ug/g	267		168	294	2	20
		Zinc		217	ug/g	311		190	352	1	20

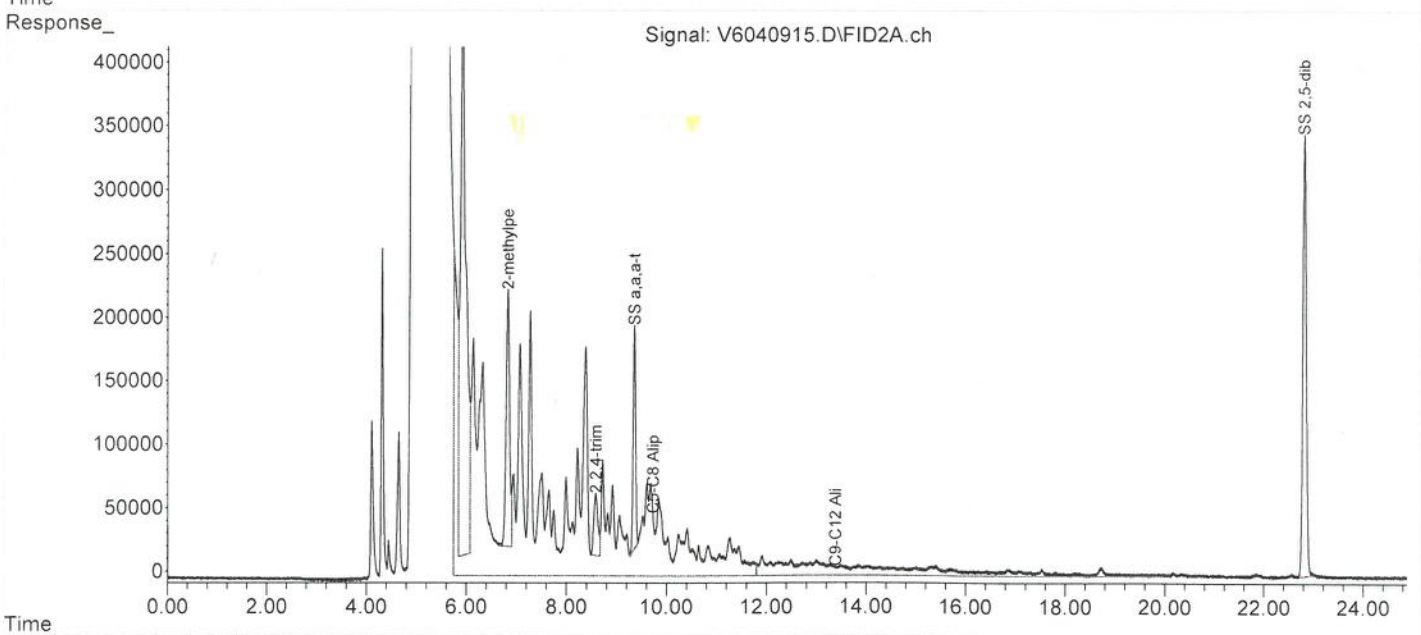
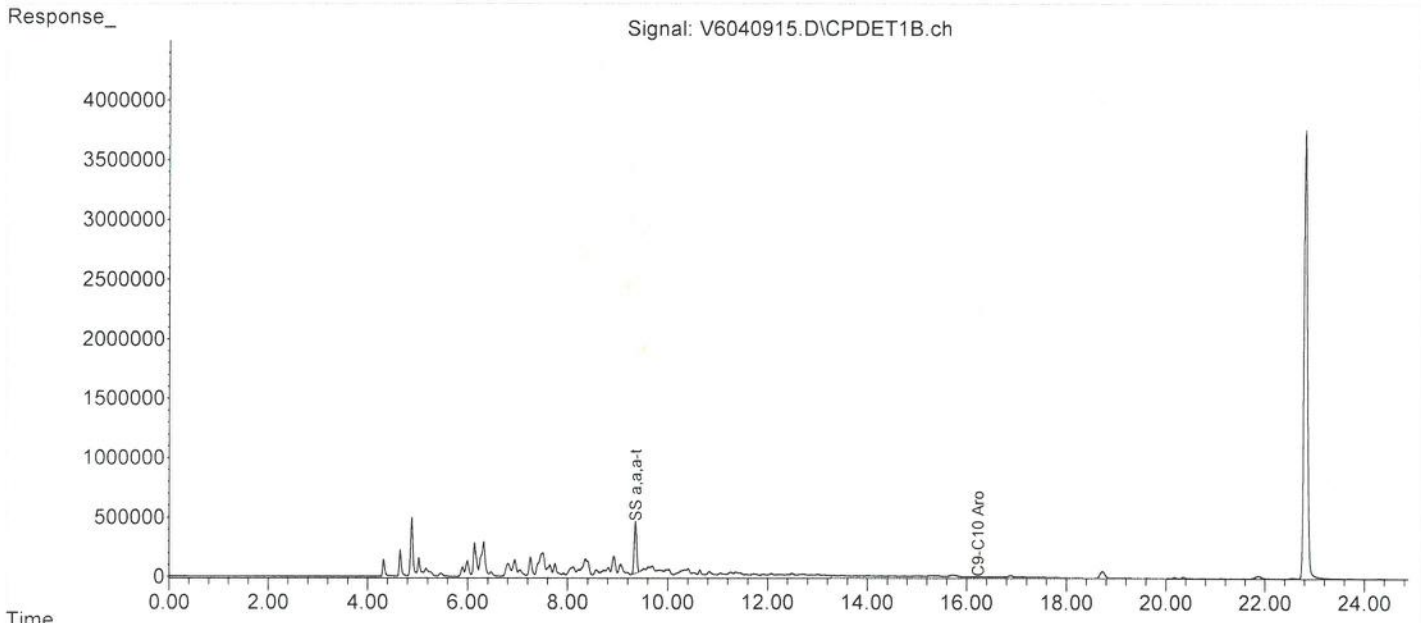
Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3051A6020A	MS12629	Silver	52513-004	170	ug/g	155	86	75	125			
		Arsenic	52513-004	280	ug/g	310	88	75	125			
		Barium	52513-004	660	ug/g	310	122	75	125			
		Beryllium	52513-004	290	ug/g	310	92	75	125			
		Cadmium	52513-004	320	ug/g	310	103	75	125			
		Chromium	52513-004	340	ug/g	310	98	75	125			
		Nickel	52513-004	340	ug/g	310	86	75	125			
		Lead	52513-004	1200	ug/g	310.2	-7	*	75	125		
		Antimony	52513-004	340	ug/g	310	93	75	125			
		Selenium	52513-004	240	ug/g	310	79	75	125			
		Thallium	52513-004	140	ug/g	155	90	75	125			
		Vanadium	52513-004	320	ug/g	310	95	75	125			
		Zinc	52513-004	2000	ug/g	310.2	-270	75	125			
SW3051A6020A	MSD12629	Silver	52513-004	160	ug/g	155	82	75	125	4	20	
		Arsenic	52513-004	290	ug/g	310	90	75	125	3	20	
		Barium	52513-004	770	ug/g	310	156	*	75	125	15	20
		Beryllium	52513-004	290	ug/g	310	94	75	125	2	20	
		Cadmium	52513-004	310	ug/g	310	99	75	125	4	20	
		Chromium	52513-004	380	ug/g	310	111	75	125	12	20	
		Nickel	52513-004	400	ug/g	310	103	75	125	15	20	
		Lead	52513-004	1300	ug/g	310.2	43	*	75	125	12	20
		Antimony	52513-004	350	ug/g	310	96	75	125	3	20	
		Selenium	52513-004	230	ug/g	310	74	*	75	125	7	20
		Thallium	52513-004	150	ug/g	155	95	75	125	5	20	
		Vanadium	52513-004	310	ug/g	310	94	75	125	1	20	
		Zinc	52513-004	2300	ug/g	310.2	-188	75	125	12	20	
SW7471B	BLK12632	Mercury		<	0.14	ug/g						
SW7471B	CRM12632	Mercury		0.226	ug/g	0.221		0.0908	0.351			
SW7471B	CRMD12632	Mercury		0.216	ug/g	0.221		0.0908	0.351	4	35	
SW7471B	MS12632	Mercury	52513-004	0.65	ug/g	0.405	12	*	80	120		
SW7471B	MSD12632	Mercury	52513-004	0.86	ug/g	0.405	64	*	80	120	28	35

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3060A7196A	CRM2001871	Chromium, Hexavalent		110	ug/g	260		93.3	350	
SW3060A7196A	CRMD2001871	Chromium, Hexavalent		100	ug/g	260		93.3	350	30
SW3060A7196A	DUP2001871	Chromium, Hexavalent	52498-013	< 0.48	ug/g					20
SW3060A7196A	MS2001871	Chromium, Hexavalent	52513-004	< 0.50	ug/g	25.01	1 *	75	125	
SW3060A7196A	MS2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.25	1 *	75	125	
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52513-004	< 0.49	ug/g	24.43	1 *	75	125	25
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.35	1 *	75	125	63
SW3060A7196A	PB2001871	Chromium, Hexavalent		< 0.40	ug/g					20
SW9045C	DUP2001840	pH	52513-002	5.6	pH					

Data Path : X:\V6\2020\Apr20\040920\  
Data File : V6040915.D  
Signal(s) : Signal #1: CPDET1B.ch Signal #2: FID2A.ch  
Acq On : 9 Apr 2020 3:49 pm  
Operator : lmm  
Sample : 52513-04 100uLMeOH/5mL  
Misc : 100  
ALS Vial : 15 Sample Multiplier: 1

Integration File signal 1: autoint1.e  
Integration File signal 2: autoint2.e  
Quant Time: Apr 09 20:07:01 2020  
Quant Method : X:\V6\methods\V6VPH040320.M  
Quant Title : VOA06 MA VPH  
QLast Update : Fri Apr 03 18:09:08 2020  
Response via : Initial Calibration  
Integrator: ChemStation

Volume Inj. :  
Signal #1 Phase : Signal #2 Phase:  
Signal #1 Info : Signal #2 Info :

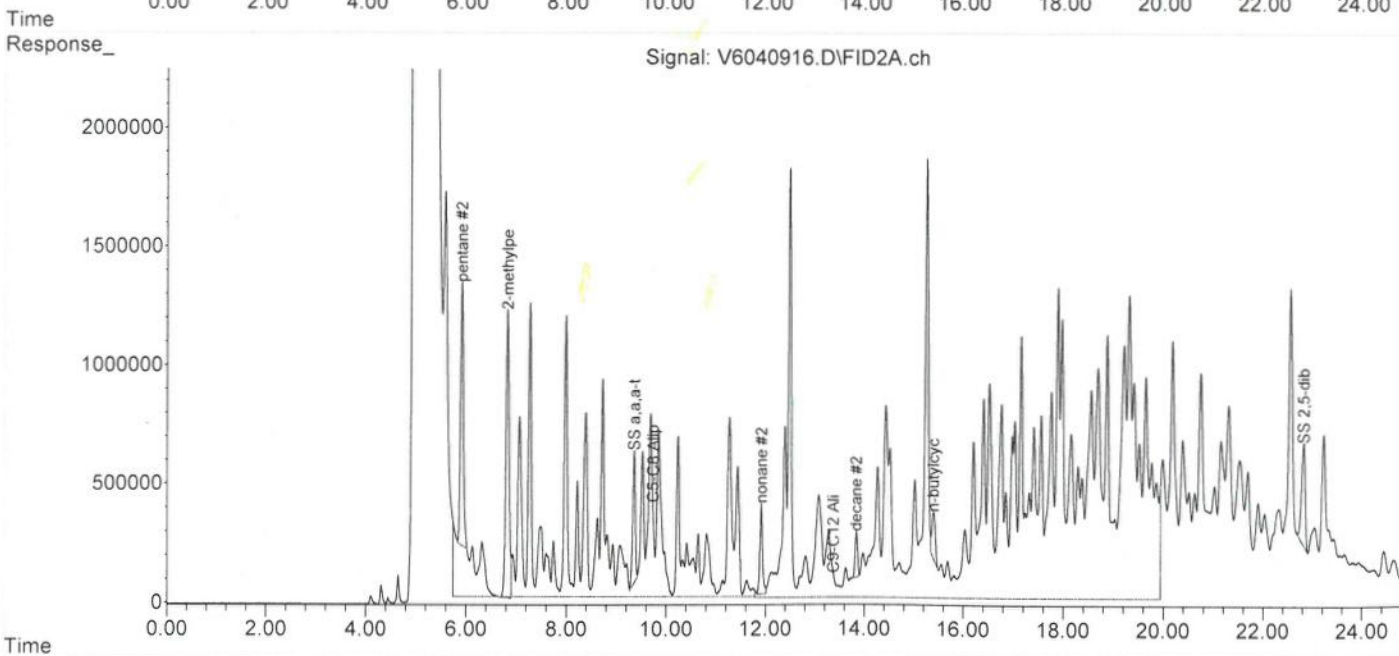
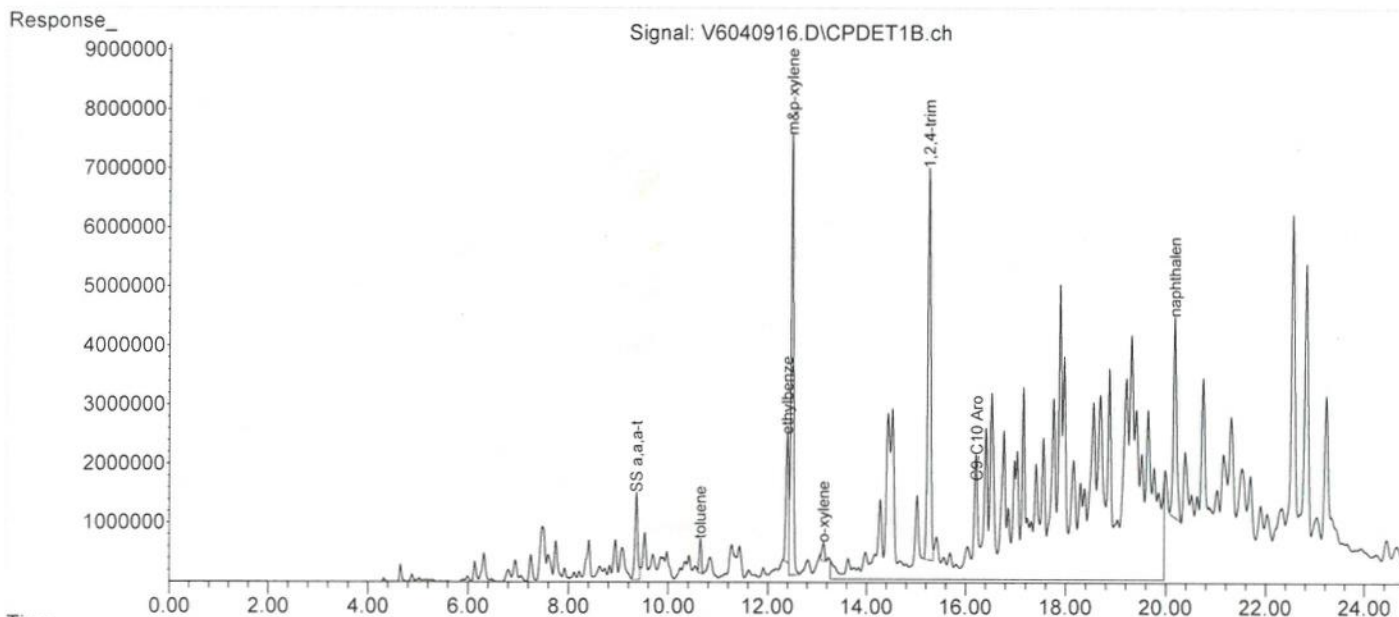




Data Path : X:\V6\2020\Apr20\040920\  
 Data File : V6040916.D  
 Signal(s) : Signal #1: CPDET1B.ch Signal #2: FID2A.ch  
 Acq On : 9 Apr 2020 4:19 pm  
 Operator : lmm  
 Sample : 52513-05 100uLMeOH/5mL  
 Misc : 100  
 ALS Vial : 16 Sample Multiplier: 1

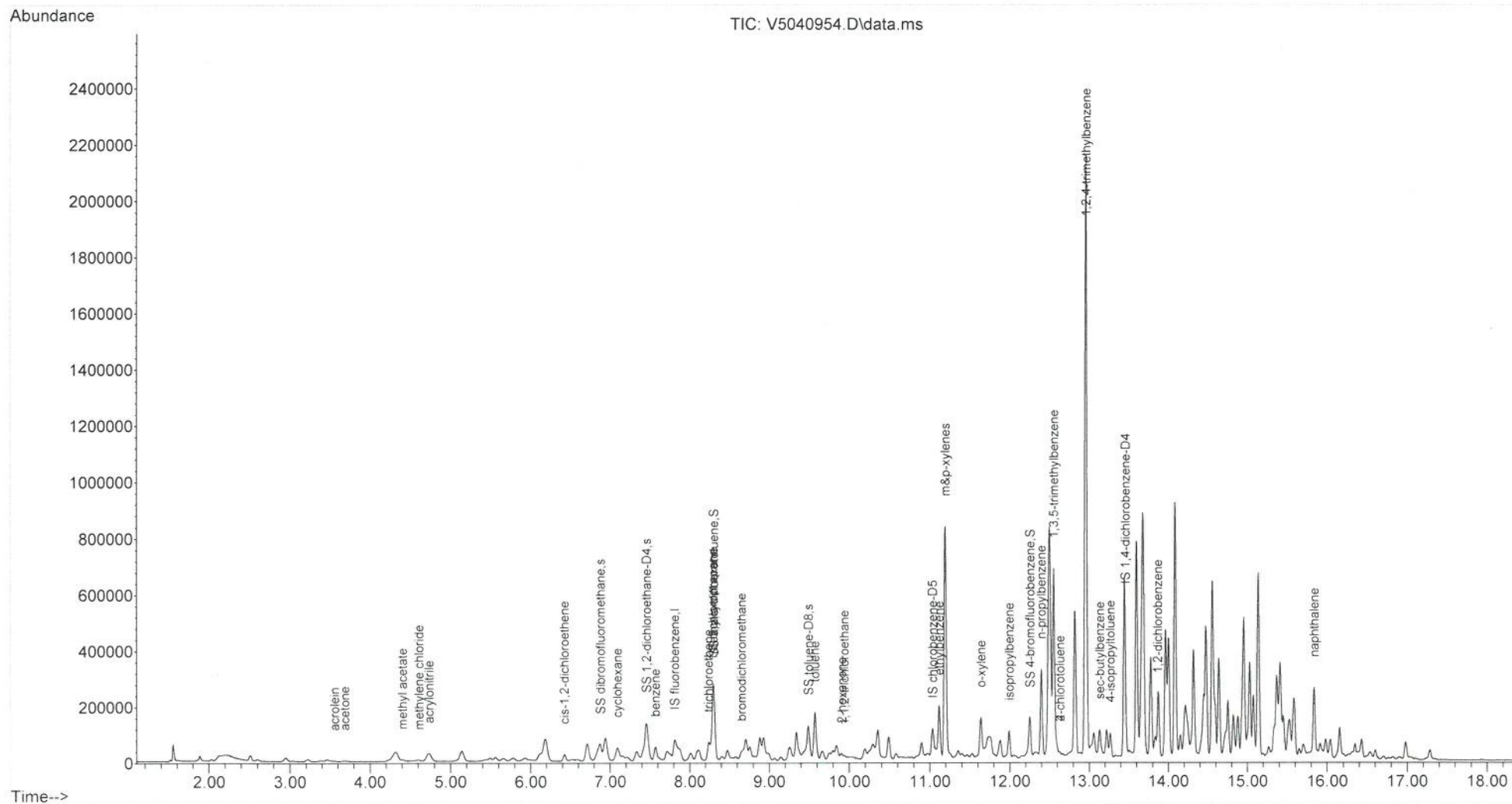
Integration File signal 1: autoint1.e  
 Integration File signal 2: autoint2.e  
 Quant Time: Apr 09 20:07:04 2020  
 Quant Method : X:\V6\methods\V6VPH040320.M  
 Quant Title : VOA06 MA VPH  
 QLast Update : Fri Apr 03 18:09:08 2020  
 Response via : Initial Calibration  
 Integrator: ChemStation

Volume Inj. :  
 Signal #1 Phase : Signal #2 Phase:  
 Signal #1 Info : Signal #2 Info :



Data Path : X:\V5\2020\Apr20\040920\  
 Data File : V5040954.D  
 Acq On : 10 Apr 2020 9:36 am  
 Operator : lmm  
 Sample : 52513-08 100uLMeOH/5mL  
 Misc : 100  
 ALS Vial : 54 Sample Multiplier: 1

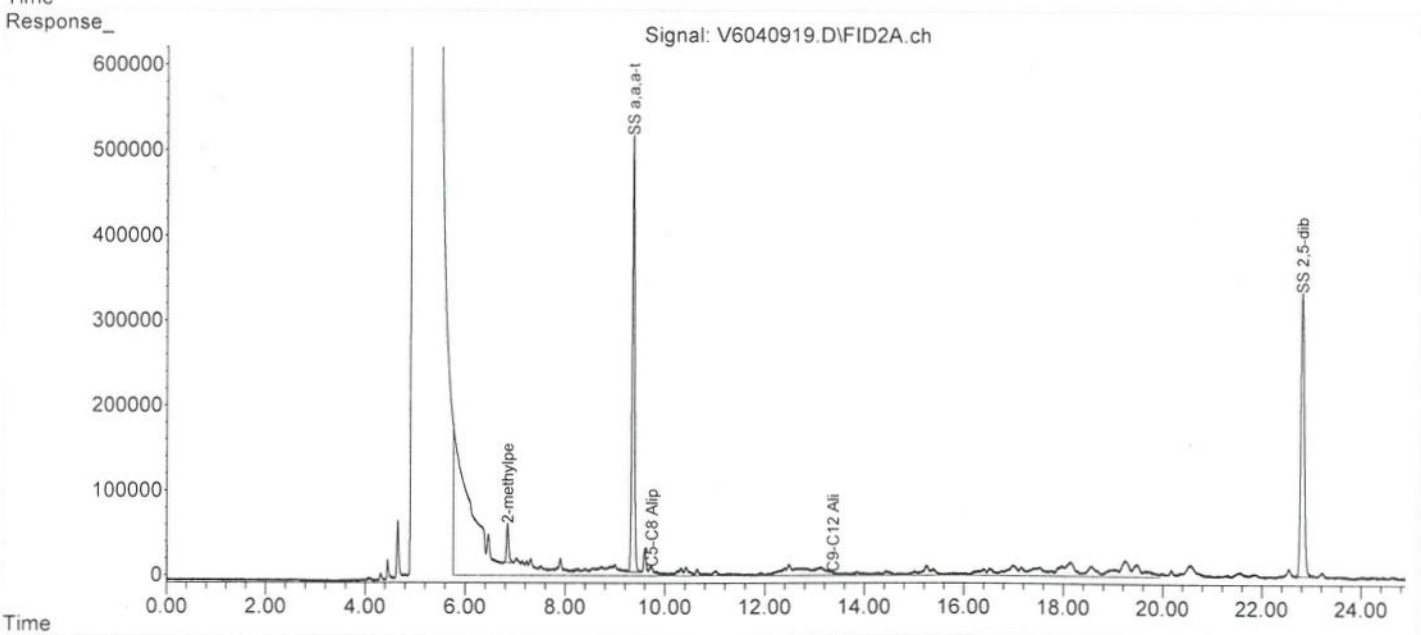
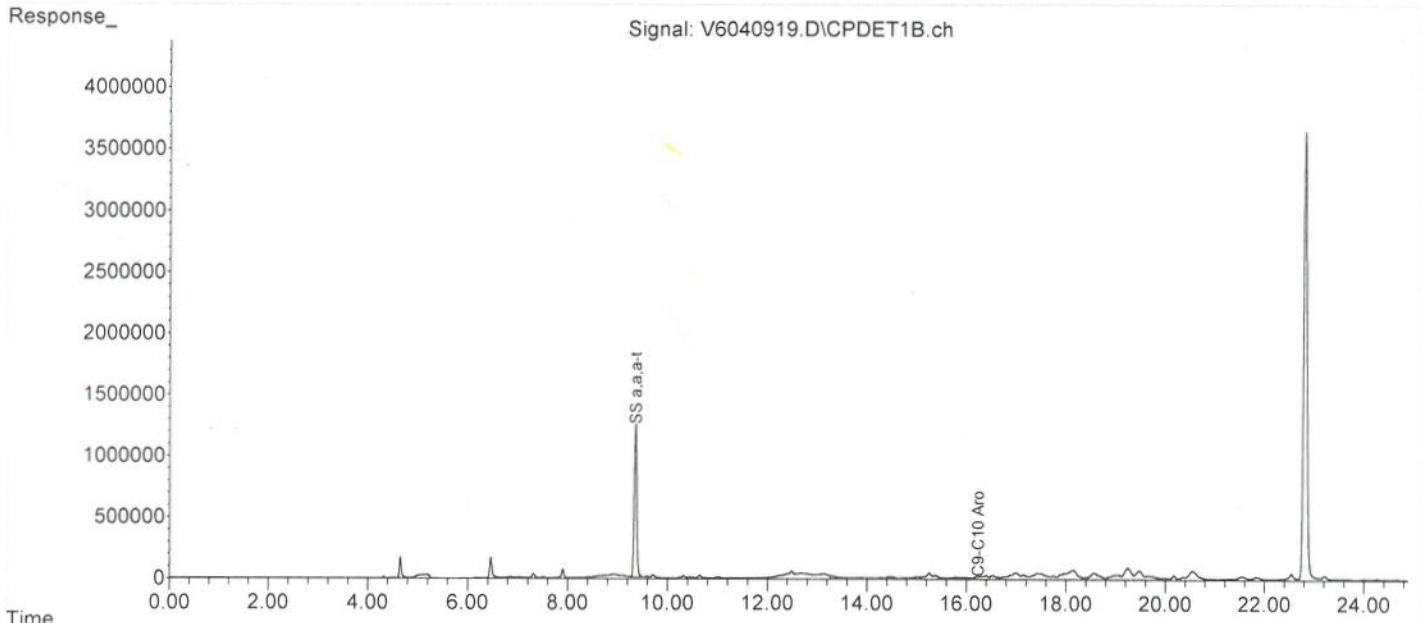
Quant Time: Apr 10 13:59:28 2020  
 Quant Method : X:\V5\methods\V58260FEB2420A.M  
 Quant Title : V5 Method 8260  
 QLast Update : Thu Apr 09 10:49:52 2020  
 Response via : Initial Calibration



Data Path : X:\V6\2020\Apr20\040920\  
Data File : V6040919.D  
Signal(s) : Signal #1: CPDET1B.ch Signal #2: FID2A.ch  
Acq On : 9 Apr 2020 5:49 pm  
Operator : lmm  
Sample : 52513-14 100uLMeOH/5mL  
Misc : 100  
ALS Vial : 19 Sample Multiplier: 1

Integration File signal 1: autoint1.e  
Integration File signal 2: autoint2.e  
Quant Time: Apr 09 20:07:13 2020  
Quant Method : X:\V6\methods\V6VPH040320.M  
Quant Title : VOA06 MA VPH  
QLast Update : Fri Apr 03 18:09:08 2020  
Response via : Initial Calibration  
Integrator: ChemStation

Volume Inj. :  
Signal #1 Phase : Signal #2 Phase:  
Signal #1 Info : Signal #2 Info :



## AROMATIC HYDROCARBON BREAKTHROUGH CALCULATION

Method: MADEP EPH 2019 Rev 2.1

	Ics12630 Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.6%	<5.0%	4/9/2020
2-methylnaphthalene	0.2%	<5.0%	4/9/2020

	Icsd12630 Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.4%	<5.0%	4/9/2020
2-methylnaphthalene	0.2%	<5.0%	4/9/2020



**Absolute Resource**  
associates



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Portsmouth, NH 03801  
603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
AND ANALYSIS REQUEST**

52513

**ANALYSIS REQUEST**

Company Name: \_\_\_\_\_  
 Company Address: SEE PAGE 1  
 Report To: \_\_\_\_\_  
 Phone #: \_\_\_\_\_  
 Invoice to: \_\_\_\_\_  
 Email: SEE PAGE 1  
 PO #: \_\_\_\_\_

Project Name: \_\_\_\_\_  
 Project #: SEE PAGE 1  
 Project Location: NH MA ME VT \_\_\_\_\_  
 Accreditation Required? N/Y: \_\_\_\_\_  
 Protocol: RCRA SDWA NPDES  
           MCP NHDES DOD  
 Reporting QAPP GW-1 S-1  
 Limits: EPA DW Other \_\_\_\_\_  
 Quote # N/A  
 NH Reimbursement Pricing

<input checked="" type="checkbox"/> VOC 8260 MADEP *	<input type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MBE only	<input type="checkbox"/> VOC 8021VT	<input checked="" type="checkbox"/> VPH MADEP *	<input type="checkbox"/> GRO 8015	<input checked="" type="checkbox"/> 1,4-Dioxane *	<input type="checkbox"/> VOC 524.2 NH List	<input type="checkbox"/> Gases-List:	<input type="checkbox"/> TPH	<input type="checkbox"/> DR0 8015	<input checked="" type="checkbox"/> EPH MADEP	<input type="checkbox"/> TPH Fingerprint	<input type="checkbox"/> 8270PAH	<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625.1	<input type="checkbox"/> EDB	<input type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608.3 Pest/PCB	<input type="checkbox"/> O&G 1664	<input type="checkbox"/> Mineral O&G 1664	<input type="checkbox"/> pH	<input type="checkbox"/> BOD	<input type="checkbox"/> Conductivity	<input type="checkbox"/> Turbidity	<input type="checkbox"/> Apparent Color	<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TS	<input type="checkbox"/> TVS	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Acidity	<input type="checkbox"/> RCRA Metals	<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input type="checkbox"/> Hardness	<input checked="" type="checkbox"/> Total Metals-list: <u>MCP 14 METALS</u>	<input type="checkbox"/> Dissolved Metals-list:	<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD	<input type="checkbox"/> TKN	<input type="checkbox"/> TN	<input type="checkbox"/> TOC	<input type="checkbox"/> Ferrrous Iron	<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Bacteria P/A	<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> Enterococci	<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide	<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Ortho P	<input type="checkbox"/> Phenols	<input type="checkbox"/> Nitrate	<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Sulfate	<input type="checkbox"/> Bromide	<input type="checkbox"/> Fluoride	<input type="checkbox"/> Corrosivity	<input type="checkbox"/> Ignitibility/FP	<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC	<input type="checkbox"/> TCLP Pesticide	<input type="checkbox"/> Subcontract:	<input type="checkbox"/> Grain Size	<input type="checkbox"/> Herbicides	<input type="checkbox"/> Asbestos	<input type="checkbox"/> PFAS	<u>CC IV</u>	Grab (G) or Composite (C)
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Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
525-312 HOLD	SB-110(0-1')			X							4/7/20	12:45	HE
13	SB-110(0-3')			X							12:55	12:55	HE
14	SB-110(5-8')			X							12:55	12:55	HE
15 HOLD	SB-112(0-1')			X							13:15	13:15	HE
16	SB-112(0-3')			X							13:20	13:20	HE
17	SB-112(5-8')			X							13:25	13:25	HE
18	MS-1			X							10:15	10:15	HE
19	MSD-1			X									HE
20	DUP-1			X				X					HE
21	TRIP BLANK							X					HE

**TAT REQUESTED**  
 Priority (24 hr)\*   
 Expedited (48 hr)\*   
 Standard (10 Business Days)   
 \*Date Needed 5/24/20

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
 \*VOCs & VPH EXCEPT FOR TRIP BLANK  
 \*\*VPH WITH TARGET VOCs

**REPORTING INSTRUCTIONS**  PDF (e-mail address) SEE PAGE 1  
 HARD COPY REQUIRED  EDD

RECEIVED ON ICE  YES  NO  
 TEMPERATURE 2 °C

**CUSTODY RECORD**  
QSD-01 Revision 11/06/19

Relinquished by Sampler:	Date: 4/7/20	Time: 1500	Received by:	Date: 4-7	Time: 3:00
Relinquished by:	Date: 4-7	Time: 16:00	Received by:	Date:	Time:
Relinquished by:	Date:	Time:	Received by Laboratory:	Date: 4/7/20	Time: 16:00

# Laboratory Report



**Absolute Resource** *associates*

124 Heritage Avenue Portsmouth NH 03801

Jill Murphy  
Weston & Sampson  
55 Walkers Brook Drive  
Reading, MA 01867

PO Number: Framingham Brownfields  
Job ID: 52537  
Date Received: 4/8/20

Project: Cedar Woods 2180311

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below. The reported results apply to the sample(s) in the condition as received at the time the laboratory took custody. This report shall not be reproduced except in full and with approval from the laboratory. The liability of ARA is limited to the cost of the requested analyses, unless otherwise agreed upon in writing.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,  
Absolute Resource Associates

A handwritten signature in black ink, appearing to read 'A. DeWees', written in a cursive style.

Aaron DeWees  
Chief Operating Officer

Date of Approval: 4/20/2020  
Total number of pages: 44

## Absolute Resource Associates Certifications

New Hampshire 1732  
Maine NH902

Massachusetts M-NH902

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-111 (0-1')	Solid	4/8/2020 9:10	52537-001	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
SB-111 (0-3')	Solid	4/8/2020 9:12	52537-002	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Field Specified Laboratory Duplicate Field Specified Matrix Spike Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020



## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-111 (4-7')	Solid	4/8/2020 9:15	52537-003	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
SB-113 (0-3')	Solid	4/8/2020 9:30	52537-004	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-113 (5-8')	Solid	4/8/2020 9:40	52537-005	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
SB-115 (0-3')	Solid	4/8/2020 10:10	52537-006	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-115 (5-8')	Solid	4/8/2020 10:20	52537-007	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
SB-114 (0-3')	Solid	4/8/2020 10:30	52537-008	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-114 (5-8')	Solid	4/8/2020 10:50	52537-009	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Hexavalent Chromium in solids by SW3060A7196A Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 ORP in solids by ASTM-1498-08 pH in solids by SW9045C Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 Zinc in solids by 6020
DUP-2	Solid	4/8/2020 0:00	52537-010	Antimony in solids by 6020 Arsenic in solids by 6020 Barium in solids by 6020 Beryllium in solids by 6020 Cadmium in solids by 6020 Chromium in solids by 6020 Combo: 14 MA MCP Metals EPH in solids by MADEP Method Lead in solids by 6020 Mercury in solids by 7471 Nickel in solids by 6020 Selenium in solids by 6020 Silver in solids by 6020 Solid Digestion for ICP Analysis Thallium in solids by 6020 Vanadium in solids by 6020 VPH in solids by MA DEP Method Zinc in solids by 6020
Trip Blank	Solid	4/8/2020 0:00	52537-011	VPH in solids by MA DEP Method

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-003

Sample ID: SB-111 (4-7')

Matrix: Solid Percent Dry: 79.9% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 9:15

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Analyst	Date	Batch	Date	Time
Unadjusted C5-C8 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
benzene	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
toluene	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
ethylbenzene	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
m&p-xylenes	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
o-xylene	< 0.11	0.11	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
naphthalene	<b>0.30</b>	0.27	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
C5-C8 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
C9-C12 Aliphatics	< 5.4	5.4	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
C9-C10 Aromatics	< 5.4	5.4	ug/g	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>107</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>108</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>67 *</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:19	MA VPH

\* **This surrogate showed recovery outside the acceptance limits.**

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-005

Sample ID: SB-113 (5-8')

Matrix: Solid Percent Dry: 87.2% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.66 mL MeOH/g soil.

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 9:40

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
benzene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
toluene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
ethylbenzene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
m&p-xylenes	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
o-xylene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
naphthalene	< 0.23	0.23	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
C5-C8 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
C9-C12 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
C9-C10 Aromatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>99</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>100</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>68 *</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	19:49	MA VPH

\* **This surrogate showed recovery outside the acceptance limits.**

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-007

Sample ID: SB-115 (5-8')

Matrix: Solid Percent Dry: 66.1% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.63 mL MeOH/g soil.

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 10:20

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 7.3	7.3	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 7.3	7.3	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
benzene	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
toluene	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
ethylbenzene	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
m&p-xylenes	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
o-xylene	< 0.15	0.15	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
naphthalene	< 0.37	0.37	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
C5-C8 Aliphatics	< 7.3	7.3	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
C9-C12 Aliphatics	< 7.3	7.3	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
C9-C10 Aromatics	< 7.3	7.3	ug/g	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>108</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>108</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>86</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-008

Sample ID: SB-114 (0-3')

Matrix: Solid Percent Dry: 93.7% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.67 mL MeOH/g soil.

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 10:30

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Analyst	Date	Batch	Date	Time
Unadjusted C5-C8 Aliphatics	< 3.9	3.9	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
Unadjusted C9-C12 Aliphatics	< 3.9	3.9	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
methyl t-butyl ether (MTBE)	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
benzene	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
toluene	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
ethylbenzene	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
m&p-xylenes	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
o-xylene	< 0.078	0.078	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
naphthalene	< 0.20	0.20	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
C5-C8 Aliphatics	< 3.9	3.9	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
C9-C12 Aliphatics	< 3.9	3.9	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
C9-C10 Aromatics	< 3.9	3.9	ug/g	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>100</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>100</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH
a,a,a-trifluorotoluene SUR	<b>79</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	20:49	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.



Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-010

Sample ID: DUP-2

Matrix: Solid Percent Dry: 85.6% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.63 mL MeOH/g soil.

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 0:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Factor	Analyst	Date	Batch	Date
Unadjusted C5-C8 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
benzene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
toluene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
ethylbenzene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
m&p-xylenes	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
o-xylene	< 0.090	0.090	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
naphthalene	< 0.23	0.23	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
C5-C8 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
C9-C12 Aliphatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
C9-C10 Aromatics	< 4.5	4.5	ug/g	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>100</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>101</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>54 *</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	21:19	MA VPH

\* **This surrogate showed recovery outside the acceptance limits.**

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-011

Sample ID: Trip Blank

Matrix: Solid

Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil

Received on ice at 1°C, in satisfactory condition.

Sampled: 4/8/20 0:00

Parameter	Reporting		Units	Instr Dil'n	Prep	Analysis				
	Result	Limit				Analyst	Date	Batch	Date	Time
Unadjusted C5-C8 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
Unadjusted C9-C12 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
methyl t-butyl ether (MTBE)	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
benzene	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
toluene	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
ethylbenzene	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
m&p-xylenes	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
o-xylene	< 0.10	0.10	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
naphthalene	< 0.25	0.25	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
C5-C8 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
C9-C12 Aliphatics	< 5.0	5.0	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
C9-C10 Aromatics	< 5.0	5.0	ug/g	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
2,5-dibromotoluene as Aromatic SUR	<b>104</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
2,5-dibromotoluene as Aliphatic SUR	<b>104</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH
a,a,a-trifluorotoluene SUR	<b>84</b>	70-130	%	1	LMM	4/9/20	12628	4/9/20	12:19	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-001

Sample ID: SB-111 (0-1')

Matrix: Solid Percent Dry: 91.8% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:10

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
2-methylnaphthalene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
phenanthrene	<b>0.57</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
acenaphthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
acenaphthylene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
fluorene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
anthracene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
fluoranthene	<b>2.2</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
pyrene	<b>1.9</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
benzo(a)anthracene	<b>1.5</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
chrysene	<b>1.8</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
benzo(b)fluoranthene	<b>2.3</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
benzo(k)fluoranthene	<b>1.7</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
benzo(a)pyrene	<b>2.0</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
indeno(1,2,3-cd)pyrene	<b>1.7</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
dibenzo(a,h)anthracene	<b>0.51</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
benzo(g,h,i)perylene	<b>2.0</b>	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	20:30	MA EPH
Unadjusted C11-C22 Aromatics	<b>80</b>	20	ug/g	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
C9-C18 Aliphatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
C19-C36 Aliphatics	<b>85</b>	20	ug/g	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
C11-C22 Aromatics	<b>62</b>	20	ug/g	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>44</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
o-terphenyl SUR	<b>48</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
2-fluorobiphenyl SUR	<b>67</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH
2-bromonaphthalene SUR	<b>61</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	8:30	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-002

Sample ID: SB-111 (0-3')

Matrix: Solid

Percent Dry: 82.2% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:12

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
phenanthrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
benzo(a)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
chrysene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
benzo(b)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
benzo(k)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
benzo(a)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
indeno(1,2,3-cd)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
dibenzo(a,h)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
benzo(g,h,i)perylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:28	MA EPH
Unadjusted C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
C9-C18 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
C19-C36 Aliphatics	<b>40</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>43</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
o-terphenyl SUR	<b>49</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
2-fluorobiphenyl SUR	<b>74</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH
2-bromonaphthalene SUR	<b>69</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:23	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-003

Sample ID: SB-111 (4-7')

Matrix: Solid Percent Dry: 79.9% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:15

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	0.35	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
2-methylnaphthalene	0.22	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
phenanthrene	5.1	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
acenaphthene	0.24	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
fluorene	0.43	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
anthracene	1.3	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
fluoranthene	8.5	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
pyrene	6.7	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
benzo(a)anthracene	4.1	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
chrysene	4.4	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
benzo(b)fluoranthene	3.9	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
benzo(k)fluoranthene	3.4	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
benzo(a)pyrene	4.1	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
indeno(1,2,3-cd)pyrene	2.7	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
dibenzo(a,h)anthracene	0.95	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
benzo(g,h,i)perylene	3.0	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	22:00	MA EPH
Unadjusted C11-C22 Aromatics	250	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
C9-C18 Aliphatics	< 97	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
C19-C36 Aliphatics	340	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
C11-C22 Aromatics	200	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	53	40-140	%	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
o-terphenyl SUR	56	40-140	%	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
2-fluorobiphenyl SUR	71	40-140	%	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH
2-bromonaphthalene SUR	62	40-140	%	5	DBV	4/15/20	12658	4/16/20	7:56	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-004

Sample ID: SB-113 (0-3')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
phenanthrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
benzo(a)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
chrysene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
benzo(b)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
benzo(k)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
benzo(a)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
indeno(1,2,3-cd)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
dibenzo(a,h)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
benzo(g,h,i)perylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	20:00	MA EPH
Unadjusted C11-C22 Aromatics	<b>99</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
C9-C18 Aliphatics	< 97	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
C19-C36 Aliphatics	<b>450</b>	97	ug/g	5	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
C11-C22 Aromatics	<b>97</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>52</b>	40-140	%	5	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
o-terphenyl SUR	<b>55</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
2-fluorobiphenyl SUR	<b>70</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	7:22	MA EPH
2-bromonaphthalene SUR	<b>64</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	7:22	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-005

Sample ID: SB-113 (5-8')

Matrix: Solid Percent Dry: 87.2% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
phenanthrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
benzo(a)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
chrysene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
benzo(b)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
benzo(k)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
benzo(a)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
indeno(1,2,3-cd)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
dibenzo(a,h)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
benzo(g,h,i)perylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	17:58	MA EPH
Unadjusted C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
C9-C18 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
C19-C36 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>52</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
o-terphenyl SUR	<b>61</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
2-fluorobiphenyl SUR	<b>72</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH
2-bromonaphthalene SUR	<b>66</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	3:57	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-006

Sample ID: SB-115 (0-3')

Matrix: Solid Percent Dry: 89.3% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:10

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
phenanthrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
benzo(a)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
chrysene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
benzo(b)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
benzo(k)fluoranthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
benzo(a)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
indeno(1,2,3-cd)pyrene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
dibenzo(a,h)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
benzo(g,h,i)perylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	18:29	MA EPH
Unadjusted C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
C9-C18 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
C19-C36 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
C11-C22 Aromatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>53</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
o-terphenyl SUR	<b>63</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
2-fluorobiphenyl SUR	<b>69</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH
2-bromonaphthalene SUR	<b>62</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:06	MA EPH



Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-007

Sample ID: SB-115 (5-8')

Matrix: Solid Percent Dry: 66.1% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	0.26	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
2-methylnaphthalene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
phenanthrene	2.2	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
acenaphthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
acenaphthylene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
fluorene	0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
anthracene	0.33	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
fluoranthene	3.3	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
pyrene	3.5	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
benzo(a)anthracene	1.6	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
chrysene	1.9	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
benzo(b)fluoranthene	1.8	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
benzo(k)fluoranthene	1.4	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
benzo(a)pyrene	1.7	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
indeno(1,2,3-cd)pyrene	1.2	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
dibenzo(a,h)anthracene	0.38	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
benzo(g,h,i)perylene	1.5	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	21:00	MA EPH
Unadjusted C11-C22 Aromatics	96	20	ug/g	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
C9-C18 Aliphatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
C19-C36 Aliphatics	57	20	ug/g	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
C11-C22 Aromatics	75	20	ug/g	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	55	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
o-terphenyl SUR	65	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
2-fluorobiphenyl SUR	71	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH
2-bromonaphthalene SUR	64	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:14	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-008

Sample ID: SB-114 (0-3')

Matrix: Solid Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
phenanthrene	<b>4.1</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
acenaphthylene	<b>0.77</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
anthracene	<b>0.61</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
fluoranthene	<b>9.3</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
pyrene	<b>11</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
benzo(a)anthracene	<b>4.9</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
chrysene	<b>5.7</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
benzo(b)fluoranthene	<b>3.6</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
benzo(k)fluoranthene	<b>3.4</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
benzo(a)pyrene	<b>4.4</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
indeno(1,2,3-cd)pyrene	<b>2.7</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
dibenzo(a,h)anthracene	<b>0.78</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
benzo(g,h,i)perylene	<b>3.4</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	21:30	MA EPH
Unadjusted C11-C22 Aromatics	<b>270</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
C9-C18 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
C19-C36 Aliphatics	<b>40</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
C11-C22 Aromatics	<b>220</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>49</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
o-terphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
2-fluorobiphenyl SUR	<b>68</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH
2-bromonaphthalene SUR	<b>63</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	5:40	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-009

Sample ID: SB-114 (5-8')

Matrix: Solid Percent Dry: 68.7% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:50

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
2-methylnaphthalene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
phenanthrene	<b>0.54</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
acenaphthene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
acenaphthylene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
fluorene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
fluoranthene	<b>0.77</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
pyrene	<b>0.69</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
benzo(a)anthracene	<b>0.33</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
chrysene	<b>0.38</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
benzo(b)fluoranthene	<b>0.29</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
benzo(k)fluoranthene	<b>0.31</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
benzo(a)pyrene	<b>0.37</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
indeno(1,2,3-cd)pyrene	<b>0.25</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
dibenzo(a,h)anthracene	< 0.19	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
benzo(g,h,i)perylene	<b>0.34</b>	0.19	ug/g	1	CL	4/15/20	12658	4/15/20	19:29	MA EPH
Unadjusted C11-C22 Aromatics	<b>38</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
C9-C18 Aliphatics	< 19	19	ug/g	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
C19-C36 Aliphatics	<b>55</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
C11-C22 Aromatics	<b>33</b>	19	ug/g	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>52</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
o-terphenyl SUR	<b>65</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
2-fluorobiphenyl SUR	<b>72</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH
2-bromonaphthalene SUR	<b>66</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	6:48	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-010

Sample ID: DUP-2

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Sampled: 4/8/20 0:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
naphthalene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
2-methylnaphthalene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
phenanthrene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
acenaphthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
acenaphthylene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
fluorene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
anthracene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
fluoranthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
pyrene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
benzo(a)anthracene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
chrysene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
benzo(b)fluoranthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
benzo(k)fluoranthene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
benzo(a)pyrene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
indeno(1,2,3-cd)pyrene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
dibenzo(a,h)anthracene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
benzo(g,h,i)perylene	< 0.20	0.20	ug/g	1	CL	4/15/20	12658	4/15/20	18:59	MA EPH
Unadjusted C11-C22 Aromatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
C9-C18 Aliphatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
C19-C36 Aliphatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
C11-C22 Aromatics	< 20	20	ug/g	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
<b>Surrogate Recovery</b>		<b>Limits</b>								
1-chloro-octadecane SUR	<b>43</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
o-terphenyl SUR	<b>49</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
2-fluorobiphenyl SUR	<b>70</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH
2-bromonaphthalene SUR	<b>63</b>	40-140	%	1	DBV	4/15/20	12658	4/16/20	4:31	MA EPH

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-001

Sample ID: SB-111 (0-1')

Matrix: Solid Percent Dry: 91.8% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:10

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	0.45	0.34	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Arsenic	< 1.7	1.7	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Barium	51	3.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Beryllium	< 0.34	0.34	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Cadmium	< 0.34	0.34	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Chromium	12	3.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Lead	21	1.7	ug/g	5	AGN	4/10/20	12634	4/14/20	23:40	SW3051A6020A
Mercury	< 0.14	0.14	ug/g	1	AGN	4/9/20	12632	4/15/20	12:50	SW7471B
Nickel	11	3.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Selenium	< 3.4	3.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Silver	< 1.7	1.7	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Thallium	< 0.34	0.34	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Vanadium	19	3.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:25	SW3051A6020A
Zinc	42	3.4	ug/g	5	AGN	4/10/20	12634	4/14/20	23:40	SW3051A6020A

Sample#: 52537-002

Sample ID: SB-111 (0-3')

Matrix: Solid Percent Dry: 82.2% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:12

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	29	0.48	ug/g	5	AGN	4/10/20	12634	4/14/20	23:50	SW3051A6020A
Arsenic	12	2.4	ug/g	5	AGN	4/10/20	12634	4/14/20	23:50	SW3051A6020A
Barium	260 M	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
M = The recovery for the matrix spike was 133%. The acceptance criteria is 75-125%.										
Beryllium	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Cadmium	3.8 DM	0.48	ug/g	5	AGN	4/10/20	12634	4/14/20	23:50	SW3051A6020A
D = The RPD for the matrix spike duplicate was outside the 20% acceptance range. M = The recovery for the matrix spike was 192%. The acceptance criteria is 75-125%.										
Chromium	43	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Lead	630 M	2.4	ug/g	5	AGN	4/10/20	12634	4/14/20	23:50	SW3051A6020A
M = The recovery for the matrix spike was 196%. The acceptance criteria is 75-125%.										
Mercury	0.26	0.18	ug/g	1	AGN	4/14/20	12647	4/16/20	16:10	SW7471B
Nickel	52	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Selenium	< 4.8 M	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
M = The recovery for the matrix spike was 72%. The acceptance criteria is 75-125%.										
Silver	4.5	2.4	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Thallium	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Vanadium	12	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	4:33	SW3051A6020A
Zinc	1500	4.8	ug/g	5	AGN	4/10/20	12634	4/14/20	23:50	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-003

Sample ID: SB-111 (4-7')

Matrix: Solid Percent Dry: 79.9% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:15

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	4.6	0.40	ug/g	5	AGN	4/10/20	12634	4/15/20	0:18	SW3051A6020A
Arsenic	8.2	2.0	ug/g	5	AGN	4/10/20	12634	4/15/20	0:18	SW3051A6020A
Barium	80	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Beryllium	0.44	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Cadmium	0.70	0.40	ug/g	5	AGN	4/10/20	12634	4/15/20	0:18	SW3051A6020A
Chromium	13	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Lead	460	2.0	ug/g	5	AGN	4/10/20	12634	4/15/20	0:18	SW3051A6020A
Mercury	< 0.16	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:52	SW7471B
Nickel	18	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Selenium	< 4.0	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Silver	< 2.0	2.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Thallium	0.40	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Vanadium	14	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	4:59	SW3051A6020A
Zinc	330	4.0	ug/g	5	AGN	4/10/20	12634	4/15/20	0:18	SW3051A6020A

Sample#: 52537-004

Sample ID: SB-113 (0-3')

Matrix: Solid Percent Dry: 81.1% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.56	0.56	ug/g	5	AGN	4/10/20	12634	4/15/20	0:55	SW3051A6020A
Arsenic	3.7	2.8	ug/g	5	AGN	4/10/20	12634	4/15/20	0:55	SW3051A6020A
Barium	48	5.6	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Beryllium	0.69	0.56	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Cadmium	< 0.56	0.56	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Chromium	15	5.6	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Lead	36	2.8	ug/g	5	AGN	4/10/20	12634	4/15/20	0:55	SW3051A6020A
Mercury	< 0.16	0.16	ug/g	1	AGN	4/9/20	12632	4/15/20	12:54	SW7471B
Nickel	< 5.6	5.6	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Selenium	< 5.6	5.6	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Silver	< 2.8	2.8	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Thallium	< 0.56	0.56	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Vanadium	16	5.6	ug/g	5	AGN	4/10/20	12634	4/11/20	5:07	SW3051A6020A
Zinc	43	5.6	ug/g	5	AGN	4/10/20	12634	4/15/20	0:55	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-005

Sample ID: SB-113 (5-8')

Matrix: Solid Percent Dry: 87.2% Results expressed on a dry weight basis.

Sampled: 4/8/20 9:40

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.42	0.42	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Arsenic	< 2.1	2.1	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Barium	<b>26</b>	4.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Beryllium	< 0.42	0.42	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Cadmium	< 0.42	0.42	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Chromium	<b>5.7</b>	4.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Lead	<b>3.2</b>	2.1	ug/g	5	AGN	4/10/20	12634	4/15/20	1:04	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/9/20	12632	4/15/20	12:56	SW7471B
Nickel	<b>6.2</b>	4.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Selenium	< 4.2	4.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Silver	< 2.1	2.1	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Thallium	< 0.42	0.42	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Vanadium	<b>7.8</b>	4.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:15	SW3051A6020A
Zinc	<b>18</b>	4.2	ug/g	5	AGN	4/10/20	12634	4/15/20	1:04	SW3051A6020A

Sample#: 52537-006

Sample ID: SB-115 (0-3')

Matrix: Solid Percent Dry: 89.3% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:10

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.40	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Arsenic	<b>2.6</b>	2.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Barium	<b>30</b>	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Beryllium	<b>0.56</b>	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Cadmium	< 0.40	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Chromium	<b>8.0</b>	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Lead	<b>14</b>	2.0	ug/g	5	AGN	4/10/20	12634	4/15/20	1:14	SW3051A6020A
Mercury	< 0.13	0.13	ug/g	1	AGN	4/9/20	12632	4/15/20	12:57	SW7471B
Nickel	<b>5.4</b>	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Selenium	< 4.0	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Silver	< 2.0	2.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Thallium	< 0.40	0.40	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Vanadium	<b>11</b>	4.0	ug/g	5	AGN	4/10/20	12634	4/11/20	5:49	SW3051A6020A
Zinc	<b>28</b>	4.0	ug/g	5	AGN	4/10/20	12634	4/15/20	1:14	SW3051A6020A

Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-007

Sample ID: SB-115 (5-8')

Matrix: Solid Percent Dry: 66.1% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	1.7	0.64	ug/g	5	AGN	4/10/20	12634	4/15/20	1:23	SW3051A6020A
Arsenic	8.3	3.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Barium	94	6.4	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Beryllium	< 0.64	0.64	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Cadmium	< 0.64	0.64	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Chromium	15	6.4	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Lead	330	3.2	ug/g	5	AGN	4/10/20	12634	4/15/20	1:23	SW3051A6020A
Mercury	< 0.21	0.21	ug/g	1	AGN	4/14/20	12647	4/16/20	16:16	SW7471B
Nickel	9.9	6.4	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Selenium	< 6.4	6.4	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Silver	< 3.2	3.2	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Thallium	< 0.64	0.64	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Vanadium	16	6.4	ug/g	5	AGN	4/10/20	12634	4/11/20	5:57	SW3051A6020A
Zinc	210	6.4	ug/g	5	AGN	4/10/20	12634	4/15/20	1:23	SW3051A6020A

Sample#: 52537-008

Sample ID: SB-114 (0-3')

Matrix: Solid Percent Dry: 93.7% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Arsenic	< 2.4	2.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Barium	17	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Beryllium	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Cadmium	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Chromium	< 4.8	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Lead	21	2.4	ug/g	5	AGN	4/10/20	12634	4/15/20	1:32	SW3051A6020A
Mercury	< 0.15	0.15	ug/g	1	AGN	4/14/20	12647	4/16/20	16:18	SW7471B
Nickel	< 4.8	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Selenium	< 4.8	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Silver	< 2.4	2.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Thallium	< 0.48	0.48	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Vanadium	< 4.8	4.8	ug/g	5	AGN	4/10/20	12634	4/11/20	6:05	SW3051A6020A
Zinc	40	4.8	ug/g	5	AGN	4/10/20	12634	4/15/20	1:32	SW3051A6020A



Project ID: Cedar Woods 2180311

Job ID: 52537

Sample#: 52537-009

Sample ID: SB-114 (5-8')

Matrix: Solid Percent Dry: 68.7% Results expressed on a dry weight basis.

Sampled: 4/8/20 10:50

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	0.66	0.59	ug/g	5	AGN	4/10/20	12634	4/15/20	1:42	SW3051A6020A
Arsenic	< 15	15	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A
Barium	210	5.9	ug/g	5	AGN	4/10/20	12634	4/11/20	6:14	SW3051A6020A
Beryllium	< 0.59	0.59	ug/g	5	AGN	4/10/20	12634	4/15/20	1:42	SW3051A6020A
Cadmium	< 0.59	0.59	ug/g	5	AGN	4/10/20	12634	4/11/20	6:14	SW3051A6020A
Chromium	< 29	29	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A
Lead	270	2.9	ug/g	5	AGN	4/10/20	12634	4/15/20	1:42	SW3051A6020A
Mercury	< 0.21	0.21	ug/g	1	AGN	4/14/20	12647	4/16/20	16:19	SW7471B
Nickel	< 29	29	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A
Selenium	< 29	29	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A
Silver	< 2.9	2.9	ug/g	5	AGN	4/10/20	12634	4/11/20	6:14	SW3051A6020A
Thallium	< 0.59	0.59	ug/g	5	AGN	4/10/20	12634	4/11/20	6:14	SW3051A6020A
Vanadium	< 29	29	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A
Zinc	340	29	ug/g	25	AGN	4/10/20	12634	4/16/20	14:46	SW3051A6020A

Sample#: 52537-010

Sample ID: DUP-2

Matrix: Solid Percent Dry: 85.6% Results expressed on a dry weight basis.

Sampled: 4/8/20 0:00

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.44	0.44	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Arsenic	< 2.2	2.2	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Barium	26	4.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Beryllium	< 0.44	0.44	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Cadmium	< 0.44	0.44	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Chromium	6.3	4.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Lead	3.4	2.2	ug/g	5	AGN	4/10/20	12634	4/15/20	1:51	SW3051A6020A
Mercury	< 0.14	0.14	ug/g	1	AGN	4/14/20	12647	4/16/20	16:21	SW7471B
Nickel	5.9	4.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Selenium	< 4.4	4.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Silver	< 2.2	2.2	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Thallium	< 0.44	0.44	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Vanadium	8.3	4.4	ug/g	5	AGN	4/10/20	12634	4/11/20	6:22	SW3051A6020A
Zinc	18	4.4	ug/g	5	AGN	4/10/20	12634	4/15/20	1:51	SW3051A6020A

**Project ID:** Cedar Woods 2180311

**Job ID:** 52537

**Sample#:** 52537-003

**Sample ID:** SB-111 (4-7')

**Matrix:** Solid Percent Dry: 79.9% Results expressed on a dry weight basis.

**Sampled:** 4/8/20 9:15

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	1.7	0.48	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	230		mV	1	WAS	2001852	4/9/20	6:15	ASTM1498-08	
pH	7.3		pH	1	WAS	2001855	4/9/20	5:57	SW9045C	

**Sample#:** 52537-005

**Sample ID:** SB-113 (5-8')

**Matrix:** Solid Percent Dry: 87.2% Results expressed on a dry weight basis.

**Sampled:** 4/8/20 9:40

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.40	0.40	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	230		mV	1	WAS	2001852	4/9/20	6:33	ASTM1498-08	
pH	8.2		pH	1	WAS	2001855	4/9/20	6:04	SW9045C	

**Sample#:** 52537-007

**Sample ID:** SB-115 (5-8')

**Matrix:** Solid Percent Dry: 66.1% Results expressed on a dry weight basis.

**Sampled:** 4/8/20 10:20

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.59	0.59	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
Oxidation Reduction Potential	160		mV	1	WAS	2001852	4/9/20	7:00	ASTM1498-08	
pH	7.0		pH	1	WAS	2001855	4/9/20	6:08	SW9045C	

**Sample#:** 52537-009

**Sample ID:** SB-114 (5-8')

**Matrix:** Solid Percent Dry: 68.7% Results expressed on a dry weight basis.

**Sampled:** 4/8/20 10:50

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Batch	Analysis		Reference
	Result	Limit	Units	Factor				Date	Time	
Chromium, Hexavalent	< 0.59 M	0.59	ug/g	1	SFM	2001871	4/10/20	11:30	SW3060A7196A	
M = The percent recovery in the matrix spike was outside acceptance criteria. See case narrative.										
Oxidation Reduction Potential	200		mV	1	WAS	2001852	4/9/20	7:08	ASTM1498-08	
pH	8.1		pH	1	WAS	2001855	4/9/20	6:15	SW9045C	

# Quality Control Report



124 Heritage Avenue Unit 16  
Portsmouth, NH 03801

[www.absoluteresourceassociates.com](http://www.absoluteresourceassociates.com)

## MassDEP Analytical Protocol Certification Form

Laboratory Name: Absolute Resource Associates

Project #: 2180311

Project Location: Massachusetts

RTN:

**This Form provides certifications for the following data set: list Laboratory Sample ID Number(s): 52537**

Matrices:  Groundwater/Surface Water  Soil/Sediment  Drinking Water  Air  Other:

**CAM Protocol** (check all that apply below):

8260 VOC CAM II A <input type="checkbox"/>	7470/7471 Hg CAM III B <input checked="" type="checkbox"/>	MassDEP VPH (GC/PID/FID) CAM IV A <input checked="" type="checkbox"/>	8082 PCB CAM V A <input type="checkbox"/>	9014 Total Cyanide/PAC CAM VI A <input type="checkbox"/>	6860 Perchlorate CAM VIII B <input type="checkbox"/>
8270 SVOC CAM II B <input type="checkbox"/>	7010 Metals CAM III C <input type="checkbox"/>	MassDEP VPH (GC/MS) CAM IV C <input type="checkbox"/>	8081 Pesticides CAM V B <input type="checkbox"/>	7196 Hex Cr CAM VI B <input checked="" type="checkbox"/>	MassDEP APH CAM IX A <input type="checkbox"/>
6010 Metals CAM III A <input type="checkbox"/>	6020 Metals CAM III D <input checked="" type="checkbox"/>	MassDEP EPH CAM IV B <input checked="" type="checkbox"/>	8151 Herbicides CAM V C <input type="checkbox"/>	8330 Explosives CAM VIII A <input type="checkbox"/>	TO-15 VOC CAM IX B <input type="checkbox"/>

**Affirmative Responses to Questions A through F are required for "Presumptive Certainty" status**

<b>A</b>	Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature) in the field or laboratory, and prepared/analyzed within method holding times?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>B</b>	Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>C</b>	Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>D</b>	Does the laboratory report comply with all the reporting requirements specified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>E</b>	VPH, EPH, APH, and TO-15 only a. VPH, EPH, and APH Methods only: Was each method conducted without significant modification(s)? (Refer to the individual method(s) for a list of significant modifications). b. APH and TO-15 Methods only: Was the complete analyte list reported for each method?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Yes <input type="checkbox"/> No
<b>F</b>	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No

**Responses to Questions G, H and I below are required for "Presumptive Certainty" status**


<b>G</b>	Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
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**Data User Note: Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40.1056 (2)(k) and WSC-07-350.**

<b>H</b>	Were all QC performance standards specified in the CAM protocol(s) achieved?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
<b>I</b>	Were results reported for the complete analyte list specified in the selected CAM protocol(s)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <sup>1</sup>

<sup>1</sup>All negative responses must be addressed in an attached laboratory narrative.

*I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, is accurate and complete.*

Signature: 

Position: Chief Operating Officer

Printed Name: Aaron DeWees

Date: 4/20/20

## Sample Integrity Table

Parameter	Method	Matrix	Minimum Volume	Recommended Container(s)	Required Preservation	Holding Time
Volatile Organics	EPA 8260	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
Volatile Organics	EPA 8260	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	14 Days
Semivolatile Organics	EPA 8270	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Semivolatile Organics	EPA 8270	Solid	20g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Organochlorine Pesticides	EPA 8081	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Organochlorine Pesticides	EPA 8081	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
PCBs	EPA 8082	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
PCBs	EPA 8082	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
Herbicides (subcontracted)	EPA 8151	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Herbicides (subcontracted)	EPA 8151	Solid	30g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
MA DEP VPH	MADEP VPH	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP VPH	MADEP VPH	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	28 Days
MA DEP EPH	MADEP EPH	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP EPH	MADEP EPH	Solid	30g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Total Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	180 Days
Dissolved Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	Filter First 1:1 HNO <sub>3</sub> to pH <2	180 Days
Total Metals	EPA 6010	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	180 Days
Total Mercury (may be combined with Total Metals)	EPA 7470	Aqueous	100mL	125mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	28 Days
Total Mercury (may be combined with Total Metals)	EPA 7471	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	28 Days
Chromium, Hexavalent	EPA 7196	Aqueous	100mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> buffer	28 Days
Chromium, Hexavalent (subcontract)	EPA 7196	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	30 Days
Cyanide, Total	EPA 9014	Aqueous	125mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ 1:1 NaOH to pH >8	14 Days
Cyanide, Total	EPA 9014	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days

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**Case Narrative**

**Lab # 52537**

**Sample Receiving and Chain of Custody Discrepancies**

Samples were received in acceptable condition, between 0 and 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

As noted on the result page, the VPH samples did not meet the 1:1 +/-25% methanol to soil ratio.

**Calibration**

SVOC: See the included table for a list of compounds quantitated by quadratic equation.

**Method Blank**

No exceptions noted.

**Surrogate Recoveries**

VPH: Samples 52537-003, -005, and -010 had a recovery for a,a,a-trifluorotoluene that was low and outside acceptance limits.

**Laboratory Control Sample Results**

VOC: The LCSD2001516 did not meet the acceptance criteria for t-butanol (TBA). Since <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

**Matrix Spike/Matrix Spike Duplicate/Duplicate Results**

Metals: The percent recovery in the MS/D of 52537-002 for barium (133%), cadmium (192%), lead (196 and 138%) and selenium (72 and 74%) were outside the acceptance criteria of 75-125%. The RPD for cadmium was 63% which was outside the acceptance limit of 20%. All other batch QC was within acceptance.

Chromium, Hexavalent: The matrix spike/duplicate for 52537-009 did not meet the acceptance criteria. The percent recovery was acceptable in the associated CRM/CRMD. The results for the ORP and pH analysis show the samples exhibits a reducing nature therefore resulting in poor spike recovery. The sample has been qualified accordingly.

**Other**

VPH: The trap used for VPH analysis is a Tekmar STRATUM Purge Trap 9. The column used for VPH analysis is a Restek Rtx-502.2, 105m, 0.53mmID, and 3um df.

**MassDEP Analytical Protocol Certification Form Questions A through I**

No explanation is needed for Questions A through I answered in the affirmative.

**Question G:** The CAM protocol reporting limits were not achieved for this project due to dilutions necessary for sample analysis. Box G is "No."

**Question H:** See surrogate section above. Box H is "No."



**Quantitation by Quadratic Equation**  
**Lab # 52537**

SVOC: Quantitation of the following compounds was based on a quadratic equation:

---

2-methylnaphthalene

benzo(k)fluoranthene

indeno(1,2,3-cd)pyrene

dibenzo(a,h)anthracene

benzo(g,h,i)perylene

## GLOSSARY

%R	Percent Recovery
BLK	Blank (Method Blank, Preparation Blank)
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRM	Certified Reference Material (associated with solid Metals samples)
CRMD	Certified Reference Material Duplicate (associated with solid Metals samples)
Dil'n	Dilution
DL	Detection Limit
DUP	Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
LOD	Limit of Detection
LOQ	Limit of Quantitation
MB	Methanol Blank (associated with solid VOC samples)
MLCS	Methanol Laboratory Control Sample (associated with solid VOC samples)
MLCSD	Methanol Laboratory Control Sample Duplicate (associated with solid VOC samples)
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PB	Preparation Blank
QC	Quality Control
RL	Reporting Limit
RPD	Relative Percent Difference
SUR	Surrogate



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- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
MA VPH	MB12628	Unadjusted C5-C8 Aliphatics	<	5.0	ug/g						
		Unadjusted C9-C12 Aliphatics	<	5.0	ug/g						
		methyl t-butyl ether (MTBE)	<	0.10	ug/g						
		benzene	<	0.10	ug/g						
		toluene	<	0.10	ug/g						
		ethylbenzene	<	0.10	ug/g						
		m&p-xylenes	<	0.10	ug/g						
		o-xylene	<	0.10	ug/g						
		naphthalene	<	0.25	ug/g						
		C5-C8 Aliphatics	<	5.0	ug/g						
		C9-C12 Aliphatics	<	5.0	ug/g						
		C9-C10 Aromatics	<	5.0	ug/g						
		2,5-dibromotoluene as Aromatic SUR		113	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		
MA VPH	MLCS12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130		
		Unadjusted C9-C12 Aliphatics		13	ug/g	15	89	70	130		
		methyl t-butyl ether (MTBE)		4.6	ug/g	5	92	70	130		
		benzene		4.8	ug/g	5	97	70	130		
		toluene		4.8	ug/g	5	95	70	130		
		ethylbenzene		4.8	ug/g	5	96	70	130		
		m&p-xylenes		9.7	ug/g	10	97	70	130		
		o-xylene		4.8	ug/g	5	96	70	130		
		naphthalene		5.0	ug/g	5	100	70	130		
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		
		C9-C10 Aromatics		5.1	ug/g	5	102	70	130		
		2,5-dibromotoluene as Aromatic SUR		111	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		112	%			70	130		
		a,a,a-trifluorotoluene SUR		86	%			70	130		
MA VPH	MLCSD12628	Unadjusted C5-C8 Aliphatics		14	ug/g	15	93	70	130	0	25
		Unadjusted C9-C12 Aliphatics		14	ug/g	15	90	70	130	2	25
		methyl t-butyl ether (MTBE)		4.7	ug/g	5	93	70	130	1	25
		benzene		4.9	ug/g	5	99	70	130	2	25
		toluene		4.9	ug/g	5	98	70	130	3	25
		ethylbenzene		4.9	ug/g	5	99	70	130	3	25
		m&p-xylenes		10.0	ug/g	10	100	70	130	3	25
		o-xylene		4.9	ug/g	5	99	70	130	2	25
		naphthalene		5.1	ug/g	5	102	70	130	2	25
		C5-C8 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C12 Aliphatics	<	5.0	ug/g			70	130		25
		C9-C10 Aromatics		5.2	ug/g	5	103	70	130	1	25
		2,5-dibromotoluene as Aromatic SUR		99	%			70	130		
		2,5-dibromotoluene as Aliphatic SUR		100	%			70	130		
		a,a,a-trifluorotoluene SUR		90	%			70	130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK12658	naphthalene		<	0.20	ug/g				
		2-methylnaphthalene		<	0.20	ug/g				
		phenanthrene		<	0.20	ug/g				
		acenaphthene		<	0.20	ug/g				
		acenaphthylene		<	0.20	ug/g				
		fluorene		<	0.20	ug/g				
		anthracene		<	0.20	ug/g				
		fluoranthene		<	0.20	ug/g				
		pyrene		<	0.20	ug/g				
		benzo(a)anthracene		<	0.20	ug/g				
		chrysene		<	0.20	ug/g				
		benzo(b)fluoranthene		<	0.20	ug/g				
		benzo(k)fluoranthene		<	0.20	ug/g				
		benzo(a)pyrene		<	0.20	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.20	ug/g				
		dibenzo(a,h)anthracene		<	0.20	ug/g				
		benzo(g,h,i)perylene		<	0.20	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			44	%		40	140	
		o-terphenyl SUR			48	%		40	140	
		2-fluorobiphenyl SUR			64	%		40	140	
		2-bromonaphthalene SUR			59	%		40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCS12658	naphthalene		2.7	ug/g	6	45	40	140	
		2-methylnaphthalene		2.8	ug/g	6	47	40	140	
		phenanthrene		3.2	ug/g	6	54	40	140	
		acenaphthene		2.6	ug/g	6	43	40	140	
		acenaphthylene		2.5	ug/g	6	42	40	140	
		fluorene		2.8	ug/g	6	46	40	140	
		anthracene		3.2	ug/g	6	53	40	140	
		fluoranthene		3.4	ug/g	6	57	40	140	
		pyrene		3.3	ug/g	6	55	40	140	
		benzo(a)anthracene		3.3	ug/g	6	54	40	140	
		chrysene		3.3	ug/g	6	54	40	140	
		benzo(b)fluoranthene		3.1	ug/g	6	51	40	140	
		benzo(k)fluoranthene		3.1	ug/g	6	51	40	140	
		benzo(a)pyrene		3.1	ug/g	6	52	40	140	
		indeno(1,2,3-cd)pyrene		3.8	ug/g	6	63	40	140	
		dibenzo(a,h)anthracene		3.8	ug/g	6	64	40	140	
		benzo(g,h,i)perylene		3.9	ug/g	6	64	40	140	
		Unadjusted C11-C22 Aromatics		60	ug/g	102	59	40	140	
		C9-C18 Aliphatics	<	20	ug/g	36	45	40	140	
		C19-C36 Aliphatics		43	ug/g	48	91	40	140	
		C11-C22 Aromatics	<	20	ug/g			40	140	
		1-chloro-octadecane SUR		47	%			40	140	
		o-terphenyl SUR		54	%			40	140	
		2-fluorobiphenyl SUR		73	%			40	140	
		2-bromonaphthalene SUR		69	%			40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD12658	naphthalene		2.9	ug/g	6	48	40 140	7	25
		2-methylnaphthalene		3.0	ug/g	6	50	40 140	7	25
		phenanthrene		3.4	ug/g	6	57	40 140	7	25
		acenaphthene		2.7	ug/g	6	45	40 140	4	25
		acenaphthylene		2.7	ug/g	6	44	40 140	6	25
		fluorene		2.9	ug/g	6	49	40 140	6	25
		anthracene		3.4	ug/g	6	56	40 140	6	25
		fluoranthene		3.6	ug/g	6	61	40 140	6	25
		pyrene		3.5	ug/g	6	59	40 140	7	25
		benzo(a)anthracene		3.6	ug/g	6	60	40 140	9	25
		chrysene		3.6	ug/g	6	59	40 140	9	25
		benzo(b)fluoranthene		3.5	ug/g	6	58	40 140	12	25
		benzo(k)fluoranthene		3.4	ug/g	6	57	40 140	10	25
		benzo(a)pyrene		3.4	ug/g	6	57	40 140	10	25
		indeno(1,2,3-cd)pyrene		4.1	ug/g	6	68	40 140	8	25
		dibenzo(a,h)anthracene		4.1	ug/g	6	69	40 140	8	25
		benzo(g,h,i)perylene		4.2	ug/g	6	70	40 140	9	25
		Unadjusted C11-C22 Aromatics		63	ug/g	102	62	40 140	4	25
		C9-C18 Aliphatics	<	20	ug/g	36	44	40 140	2	25
		C19-C36 Aliphatics		44	ug/g	48	91	40 140	0	25
		C11-C22 Aromatics	<	20	ug/g			40 140		25
		1-chloro-octadecane SUR		46	%			40 140		
		o-terphenyl SUR		54	%			40 140		
		2-fluorobiphenyl SUR		69	%			40 140		
		2-bromonaphthalene SUR		64	%			40 140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3051A6020A	BLK12634	Silver		<	2.5	ug/g					
		Arsenic		<	2.5	ug/g					
		Barium		<	5.0	ug/g					
		Beryllium		<	0.50	ug/g					
		Cadmium		<	0.50	ug/g					
		Chromium		<	5.0	ug/g					
		Nickel		<	5.0	ug/g					
		Lead		<	2.5	ug/g					
		Antimony		<	0.50	ug/g					
		Selenium		<	5.0	ug/g					
		Thallium		<	0.50	ug/g					
		Vanadium		<	5.0	ug/g					
		Zinc		<	5.0	ug/g					
SW3051A6020A	CRM12634	Silver		44.4	ug/g	53.8		31.9	59.7		
		Arsenic		169	ug/g	219		129	240		
		Barium		582	ug/g	788		509	867		
		Beryllium		201	ug/g	247		160	272		
		Cadmium		145	ug/g	175		111	192		
		Chromium		286	ug/g	375		223	414		
		Nickel		248	ug/g	318		193	358		
		Lead		239	ug/g	321		207	353		
		Antimony		81.0	ug/g	159		15.9	180		
		Selenium		102	ug/g	145		78.5	160		
		Thallium		81.9	ug/g	98.6		53.7	111		
		Vanadium		186	ug/g	267		168	294		
		Zinc		234	ug/g	311		190	352		
SW3051A6020A	CRMD12634	Silver		46.5	ug/g	53.8		31.9	59.7	4	20
		Arsenic		178	ug/g	219		129	240	5	20
		Barium		597	ug/g	788		509	867	3	20
		Beryllium		205	ug/g	247		160	272	2	20
		Cadmium		151	ug/g	175		111	192	4	20
		Chromium		291	ug/g	375		223	414	2	20
		Nickel		251	ug/g	318		193	358	1	20
		Lead		245	ug/g	321		207	353	3	20
		Antimony		82.8	ug/g	159		15.9	180	2	20
		Selenium		106	ug/g	145		78.5	160	4	20
		Thallium		85.1	ug/g	98.6		53.7	111	4	20
		Vanadium		188	ug/g	267		168	294	1	20
		Zinc		237	ug/g	311		190	352	1	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit		
SW3051A6020A	MS12634	Silver	52537-002	130	ug/g	133	97	75	125			
		Arsenic	52537-002	240	ug/g	266	87	75	125			
		Barium	52537-002	550	ug/g	266	109	75	125			
		Beryllium	52537-002	260	ug/g	266	99	75	125			
		Cadmium	52537-002	280	ug/g	266	103	75	125			
		Chromium	52537-002	290	ug/g	266	91	75	125			
		Nickel	52537-002	270	ug/g	266	83	75	125			
		Lead	52537-002	1200	ug/g	266.7	196	*	75	125		
		Antimony	52537-002	310	ug/g	266	105		75	125		
		Selenium	52537-002	190	ug/g	266	72	*	75	125		
		Thallium	52537-002	120	ug/g	133	89		75	125		
		Vanadium	52537-002	250	ug/g	266	91		75	125		
		Zinc	52537-002	1800	ug/g	266.7	107		75	125		
SW3051A6020A	MSD12634	Silver	52537-002	140	ug/g	138	98	75	125	5	20	
		Arsenic	52537-002	250	ug/g	276	87	75	125	3	20	
		Barium	52537-002	630	ug/g	276	133	*	75	125	13	20
		Beryllium	52537-002	280	ug/g	276	101		75	125	6	20
		Cadmium	52537-002	540	ug/g	276	192	*	75	125	63	* 20
		Chromium	52537-002	300	ug/g	276	92		75	125	5	20
		Nickel	52537-002	300	ug/g	276	91		75	125	10	20
		Lead	52537-002	1000	ug/g	276.4	138	*	75	125	13	20
		Antimony	52537-002	320	ug/g	276	106		75	125	4	20
		Selenium	52537-002	200	ug/g	276	74	*	75	125	6	20
		Thallium	52537-002	130	ug/g	138	93		75	125	8	20
		Vanadium	52537-002	260	ug/g	276	91		75	125	3	20
		Zinc	52537-002	2000	ug/g	276.4	187		75	125	12	20
SW7471B	BLK12632	Mercury		<	0.14	ug/g						
SW7471B	CRM12632	Mercury		0.226	ug/g	0.221		0.0908	0.351			
SW7471B	CRMD12632	Mercury		0.216	ug/g	0.221		0.0908	0.351	4	35	
SW7471B	MS12632	Mercury	52513-004	0.65	ug/g	0.405	12	*	80	120		
SW7471B	MSD12632	Mercury	52513-004	0.86	ug/g	0.405	64	*	80	120	28	35
SW7471B	BLK12647	Mercury		<	0.14	ug/g						
SW7471B	CRM12647	Mercury		0.198	ug/g	0.221		0.0908	0.351			
SW7471B	CRMD12647	Mercury		0.203	ug/g	0.221		0.0908	0.351	2	35	
SW7471B	MS12647	Mercury	52537-002	0.58	ug/g	0.389	82		80	120		
SW7471B	MSD12647	Mercury	52537-002	0.61	ug/g	0.374	95		80	120	6	35

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3060A7196A	CRM2001871	Chromium, Hexavalent		110	ug/g	260		93.3	350	
SW3060A7196A	CRMD2001871	Chromium, Hexavalent		100	ug/g	260		93.3	350	30
SW3060A7196A	DUP2001871	Chromium, Hexavalent	52498-013	< 0.48	ug/g					20
SW3060A7196A	MS2001871	Chromium, Hexavalent	52513-004	< 0.50	ug/g	25.01	1 *	75	125	
SW3060A7196A	MS2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.25	1 *	75	125	
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52513-004	< 0.49	ug/g	24.43	1 *	75	125	25
SW3060A7196A	MSD2001871	Chromium, Hexavalent	52537-009	< 0.55	ug/g	27.35	1 *	75	125	63
SW3060A7196A	PB2001871	Chromium, Hexavalent		< 0.40	ug/g					
SW9045C	DUP2001855	pH	52537-003	7.2	pH					

## AROMATIC HYDROCARBON BREAKTHROUGH CALCULATION

Method: MADEP EPH 2019 Rev 2.1

	Ics12658	Acceptance	Date of Analysis
	Aliphatic Breakthrough	Criteria	
	(%)		
naphthalene	0.6%	<5.0%	4/16/2020
2-methylnaphthalene	0.1%	<5.0%	4/16/2020

	Icsd12658	Acceptance	Date of Analysis
	Aliphatic Breakthrough	Criteria	
	(%)		
naphthalene	0.9%	<5.0%	4/16/2020
2-methylnaphthalene	0.1%	<5.0%	4/16/2020





124 Heritage Avenue #16  
Portsmouth, NH 03801  
603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
AND ANALYSIS REQUEST**

52537

**ANALYSIS REQUEST**

Company Name: **WESTON & SAMPSON**  
Company Address: **55 WALKERS BROOK DR READING, MA**  
Report To: **SARAH DESTEFANO & JILL MURPHY**  
Phone #: **1 800 SAMPSON**  
Invoice to: **SARAH DESTEFANO**  
Email: **DESTEFANOS@WSEINC.COM**  
PO #: **FRAMINGHAM BROWNFIELDS**

Project Name: **CEDAR WOODS**  
Project #: **218031**  
Project Location: **NH MA ME VT**  
Accreditation Required? **N/Y:**  
Protocol: **RCRA SDWA NPDES**  
**MCP NHDES DOD**  
Reporting: **QAPP GW-1 S-1**  
Limits: **EPA DW Other**  
Quote #: **N/A**  
 NH Reimbursement Pricing

<input type="checkbox"/> VOC 8280	<input type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MADEP	<input type="checkbox"/> VOC 624.1	<input type="checkbox"/> VOC BTEX MIBE, only	<input type="checkbox"/> VOC 8021VT	<input checked="" type="checkbox"/> VPH MADEP	<input type="checkbox"/> GRO 8015	<input type="checkbox"/> 1,4-Dioxane *	<input type="checkbox"/> VOC 524.2	<input type="checkbox"/> NH List	<input type="checkbox"/> Gases-List:	<input type="checkbox"/> TPH	<input type="checkbox"/> DR0 8015	<input checked="" type="checkbox"/> LEPH MADEP	<input type="checkbox"/> TPH Fingerprint	<input type="checkbox"/> 8270PAH	<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625.1	<input type="checkbox"/> EDB	<input type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608.3 Pest/PCB	<input type="checkbox"/> O&G 1664	<input type="checkbox"/> Mineral O&G 1664	<input type="checkbox"/> pH	<input type="checkbox"/> BOD	<input type="checkbox"/> Conductivity	<input type="checkbox"/> Turbidity	<input type="checkbox"/> Apparent Color	<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TS	<input type="checkbox"/> TVS	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Acidity	<input type="checkbox"/> RCRA Metals	<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input type="checkbox"/> Hardness	<input checked="" type="checkbox"/> Total Metals-list: <b>MCP 14 METALS</b>	<input type="checkbox"/> Dissolved Metals-list:	<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD	<input type="checkbox"/> TKN	<input type="checkbox"/> TN	<input type="checkbox"/> TOC	<input type="checkbox"/> Ferrrous Iron	<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Bacteria P/A	<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> Enterococci	<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide	<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Ortho P	<input type="checkbox"/> Phenols	<input type="checkbox"/> Nitrate	<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Sulfate	<input type="checkbox"/> Bromide	<input type="checkbox"/> Fluoride	<input type="checkbox"/> Corrosivity	<input type="checkbox"/> Ignitibility/FP	<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC	<input type="checkbox"/> TCLP Pesticide	<input type="checkbox"/> Subcontract: <input type="checkbox"/> Grain Size	<input type="checkbox"/> Herbicides	<input type="checkbox"/> Asbestos	<input type="checkbox"/> PFAS	<b>GR-IV-4/8/20 CR-VI</b>	Grab (G) or Composite (C)
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Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
52537-01	SB-111(0-1')			X							4/8/20	0910	Mix
-02	SB-111(0-3')			X								0912	Mix
-03	SB-111(4-7')			X					X			0915	
-04	SB-113(0-3')			X								0935	
-05	SB-113(5-8')			X					X			0940	
-06	SB-115(0-3')			X								1010	
-07	SB-115(5-8')			X					X			1020	
-08	SB-114(0-3')			X					X			1030	
-09	SB-114(5-8')			X								0950	
	MS-2**			X								0917	
	MSD-2**			X								0917	

**TAT REQUESTED**  
Priority (24 hr)\*   
Expedited (48 hr)\*   
Standard   
(10 Business Days)  
\*Date Needed: **5 DAY**

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
\* VPH w/ TARGET VOCs (0-3) per Jill 4/9/20 g  
\*\* ASSOCIATED w/ SAMPLE SB-111(2-5)

**REPORTING INSTRUCTIONS**  PDF (e-mail address) **MURPHY J@WSEINC.COM**  
 HARD COPY REQUIRED  EDD

RECEIVED ON ICE  YES  NO  
TEMPERATURE **1** °C

**CUSTODY RECORD**  
QSD-01 Revision 11/06/19

Relinquished by Sampler: <i>Margaret Ann</i>	Date 4/8/2020	Time 1300	Received by: <i>[Signature]</i>	Date 4-8	Time 1 PM
Relinquished by: <i>[Signature]</i>	Date 4-8	Time 16:32	Received by: <i>[Signature]</i>	Date	Time
Relinquished by:	Date	Time	Received by Laboratory: <i>[Signature]</i>	Date 4/8/20	Time 16:32



124 Heritage Avenue #16  
 Portsmouth, NH 03801  
 603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
 AND ANALYSIS REQUEST**

52537

**ANALYSIS REQUEST**

Company Name: \_\_\_\_\_  
 Company Address: \_\_\_\_\_  
 Report To: SEE PAGE 1  
 Phone #: \_\_\_\_\_  
 Invoice to: \_\_\_\_\_  
 Email: \_\_\_\_\_  
 PO #: \_\_\_\_\_

Project Name: \_\_\_\_\_  
 Project #: SEE PAGE 1  
 Project Location: NH (MA) ME VT \_\_\_\_\_  
 Accreditation Required? N/Y: \_\_\_\_\_  
 Protocol: RCBA SDWA NPDES  
 (MCP) NHDES DOD  
 Reporting QAPP GW-1 (S-1)  
 Limits: EPA DW Other \_\_\_\_\_  
 Quote # N/A  
 NH Reimbursement Pricing

VOC 8260 MADEP  VOC 8260 NHDES  VOC 8260 MADEP  
 VOC 624.1  VOC BTEX MIBE only  VOC 8021VT  
 VPH MADEP  GRO 8015  1,4-Dioxane  
 VOC 524.2  VOC 524.2 NH List  Gases-List:  
 TPH  DR0 8015  EPH MADEP  TPH Fingerprint  
 8270PAH  8270ABN  625.1  EDB  
 8082 PCB  8081 Pesticides  608.3 Pest/PCB  
 O&G 1664  Mineral O&G 1664  
 pH  BOD  Conductivity  Turbidity  Apparent Color  
 TSS  TDS  TS  TVS  Alkalinity  Acidity  
 RCRA Metals  Priority Pollutant Metals  TAL Metals  Hardness  
 Total Metals-list: MCP 14 METALS  
 Dissolved Metals-list:  
 Ammonia  COD  TKN  TON  TOC  Ferrous Iron  
 T-Phosphorus  Bacteria P/A  Bacteria MPN  Enterococci  
 Cyanide  Sulfide  Nitrate + Nitrite  Ortho P  Phenols  
 Nitrate  Nitrite  Chloride  Sulfate  Bromide  Fluoride  
 Corrosivity  Ignitibility/FP  
 TCLP Metals  TCLP VOC  TCLP SVOC  TCLP Pesticide  
 Subcontract:  Grain Size  Herbicides  Asbestos  PFAS  
 CR-11 MDM 4/8/20 CR VII

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
5253740	DUP-2			X						X	4/8/20		MDM
-11	TRIP BLANK		X								4/8/20		X
											4/8/20		

**TAT REQUESTED**  
 Priority (24 hr)\*   
 Expedited (48 hr)\*   
 Standard (10 Business Days)   
 \*Date Needed: 5/14

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.  
**REPORTING INSTRUCTIONS**  
 PDF (e-mail address) SEE PAGE 1  
 HARD COPY REQUIRED  EDD

**SPECIAL INSTRUCTIONS**  
 \*VPH w/ TARGET VOCs

RECEIVED ON ICE  YES  NO  
 TEMPERATURE 1 °C

**CUSTODY RECORD**  
 QSD-01 Revision 11/06/19

Relinquished by Sampler: Margie Ann	Date: 4/8/2020 Time: 1300	Received by: <i>[Signature]</i>	Date: 4-8 Time: 10 AM
Relinquished by: <i>[Signature]</i>	Date: 4-8 Time: 16:32	Received by: <i>[Signature]</i>	Date: _____ Time: _____
Relinquished by: <i>[Signature]</i>	Date: _____ Time: _____	Received by Laboratory: <i>[Signature]</i>	Date: 4/8/20 Time: 16:52

# Laboratory Report



**Absolute Resource** *associates*

124 Heritage Avenue Portsmouth NH 03801

Jill Murphy  
Weston & Sampson  
55 Walkers Brook Drive  
Reading, MA 01867

PO Number: Framingham Brownfields  
Job ID: 53042 & 53078  
Date Received: 5/15/20

Project: Cedar Woods 21

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below. The reported results apply to the sample(s) in the condition as received at the time the laboratory took custody. This report shall not be reproduced except in full and with approval from the laboratory. The liability of ARA is limited to the cost of the requested analyses, unless otherwise agreed upon in writing.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,  
Absolute Resource Associates

A handwritten signature in black ink, appearing to read 'A. DeWees', written in a cursive style.

Aaron DeWees  
Chief Operating Officer

Date of Approval: 6/1/2020  
Total number of pages: 14

## Absolute Resource Associates Certifications

New Hampshire 1732  
Maine NH902

Massachusetts M-NH902

## Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-107 (0-1')	Solid	4/7/2020 9:45	53042-001	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-108 (0-1')	Solid	4/7/2020 10:45	53042-002	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-109 (0-1')	Solid	4/7/2020 11:30	53042-003	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-110 (0-1')	Solid	4/7/2020 12:45	53042-004	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-112 (0-1')	Solid	4/7/2020 13:15	53042-005	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-106 (0-1')	Solid	4/6/2020 9:20	53078-001	Lead in solids by 6020 Solids Digestion for ICPMS Analysis
SB-102 (0-1')	Solid	4/6/2020 11:30	53078-002	Lead in solids by 6020 Solids Digestion for ICPMS Analysis

**Project ID:** Cedar Woods 21

**Job ID:** 53042

**Sample#:** 53042-001

**Sample ID:** SB-107 (0-1')

**Matrix:** Solid Percent Dry: 87.3% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 9:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>67</b>	2.7	ug/g	5	EEB	5/26/20	12797	5/28/20	19:41	SW3051A6020A

**Sample#:** 53042-002

**Sample ID:** SB-108 (0-1')

**Matrix:** Solid Percent Dry: 75.8% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 10:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>900</b>	3.1	ug/g	5	EEB	5/26/20	12797	5/28/20	19:49	SW3051A6020A

**Sample#:** 53042-003

**Sample ID:** SB-109 (0-1')

**Matrix:** Solid Percent Dry: 81.6% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 11:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>790</b>	3.0	ug/g	5	EEB	5/26/20	12797	5/28/20	19:57	SW3051A6020A

**Sample#:** 53042-004

**Sample ID:** SB-110 (0-1')

**Matrix:** Solid Percent Dry: 91.6% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 12:45

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>11</b>	2.7	ug/g	5	EEB	5/26/20	12797	5/28/20	20:06	SW3051A6020A

**Sample#:** 53042-005

**Sample ID:** SB-112 (0-1')

**Matrix:** Solid Percent Dry: 93.3% Results expressed on a dry weight basis.

**Sampled:** 4/7/20 13:15

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>25</b>	2.6	ug/g	5	EEB	5/26/20	12797	5/28/20	20:14	SW3051A6020A

**Project ID:** Cedar Woods 2180311

**Job ID:** 53078

**Sample#:** 53078-001

**Sample ID:** SB-106 (0-1')

**Matrix:** Solid      Percent Dry: 87.9% Results expressed on a dry weight basis.

**Sampled:** 4/6/20    9:20

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>68</b>	2.3	ug/g	5	EEB	5/26/20	12797	5/29/20	18:14	SW3051A6020A

**Sample#:** 53078-002

**Sample ID:** SB-102 (0-1')

**Matrix:** Solid      Percent Dry: 94.4% Results expressed on a dry weight basis.

**Sampled:** 4/6/20    11:30

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Lead	<b>73</b>	2.5	ug/g	5	EEB	5/26/20	12797	5/29/20	18:22	SW3051A6020A

# Quality Control Report



124 Heritage Avenue Unit 16  
Portsmouth, NH 03801  
[www.absoluteresourceassociates.com](http://www.absoluteresourceassociates.com)

## MassDEP Analytical Protocol Certification Form

Laboratory Name: Absolute Resource Associates

Project #: 2180311

Project Location: Massachusetts

RTN:

**This Form provides certifications for the following data set: list Laboratory Sample ID Number(s): 53042&53078**

Matrices:  Groundwater/Surface Water  Soil/Sediment  Drinking Water  Air  Other:

**CAM Protocol** (check all that apply below):

8260 VOC CAM II A <input type="checkbox"/>	7470/7471 Hg CAM III B <input type="checkbox"/>	MassDEP VPH (GC/PID/FID) CAM IV A <input type="checkbox"/>	8082 PCB CAM V A <input type="checkbox"/>	9014 Total Cyanide/PAC CAM VI A <input type="checkbox"/>	6860 Perchlorate CAM VIII B <input type="checkbox"/>
8270 SVOC CAM II B <input type="checkbox"/>	7010 Metals CAM III C <input type="checkbox"/>	MassDEP VPH (GC/MS) CAM IV C <input type="checkbox"/>	8081 Pesticides CAM V B <input type="checkbox"/>	7196 Hex Cr CAM VI B <input type="checkbox"/>	MassDEP APH CAM IX A <input type="checkbox"/>
6010 Metals CAM III A <input type="checkbox"/>	6020 Metals CAM III D <input checked="" type="checkbox"/>	MassDEP EPH CAM IV B <input type="checkbox"/>	8151 Herbicides CAM V C <input type="checkbox"/>	8330 Explosives CAM VIII A <input type="checkbox"/>	TO-15 VOC CAM IX B <input type="checkbox"/>

**Affirmative Responses to Questions A through F are required for "Presumptive Certainty" status**

<b>A</b>	Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature) in the field or laboratory, and prepared/analyzed within method holding times?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>B</b>	Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>C</b>	Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>D</b>	Does the laboratory report comply with all the reporting requirements specified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
<b>E</b>	VPH, EPH, APH, and TO-15 only a. VPH, EPH, and APH Methods only: Was each method conducted without significant modification(s)? (Refer to the individual method(s) for a list of significant modifications). b. APH and TO-15 Methods only: Was the complete analyte list reported for each method?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Yes <input type="checkbox"/> No
<b>F</b>	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No

**Responses to Questions G, H and I below are required for "Presumptive Certainty" status**


<b>G</b>	Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
----------	---	--

**Data User Note:** Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40. 1056 (2)(k) and WSC-07-350.

<b>H</b>	Were all QC performance standards specified in the CAM protocol(s) achieved?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>
<b>I</b>	Were results reported for the complete analyte list specified in the selected CAM protocol(s)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <sup>1</sup>

<sup>1</sup>All negative responses must be addressed in an attached laboratory narrative.

*I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, is accurate and complete.*

Signature: 

Position: Chief Operating Officer

Printed Name: Aaron DeWees

Date: 6/1/20



## Sample Integrity Table

Parameter	Method	Matrix	Minimum Volume	Recommended Container(s)	Required Preservation	Holding Time
Volatile Organics	EPA 8260	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
Volatile Organics	EPA 8260	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	14 Days
Semivolatile Organics	EPA 8270	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Semivolatile Organics	EPA 8270	Solid	20g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Organochlorine Pesticides	EPA 8081	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Organochlorine Pesticides	EPA 8081	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
PCBs	EPA 8082	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
PCBs	EPA 8082	Solid	20g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	365 Days
Herbicides (subcontracted)	EPA 8151	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	7 Days
Herbicides (subcontracted)	EPA 8151	Solid	30g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
MA DEP VPH	MADEP VPH	Aqueous	40mL	2 x 40mL VOA Vials with Teflon lined septa	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP VPH	MADEP VPH	Solid	40mL	1 x 40mL VOA Vial with 10mLs Methanol <u>and</u> 1 unpreserved container for percent moisture	Cool to $\leq 6^{\circ}\text{C}$ Methanol	28 Days
MA DEP EPH	MADEP EPH	Aqueous	1L	1L Amber Glass Bottle w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$ 1:1 HCl to pH <2	14 Days
MA DEP EPH	MADEP EPH	Solid	30g	4oz Amber Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days
Total Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	180 Days
Dissolved Metals	EPA 6010	Aqueous	100mL	250mL Polyethylene Bottle	Filter First 1:1 HNO <sub>3</sub> to pH <2	180 Days
Total Metals	EPA 6010	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	180 Days
Total Mercury (may be combined with Total Metals)	EPA 7470	Aqueous	100mL	125mL Polyethylene Bottle	1:1 HNO <sub>3</sub> to pH <2	28 Days
Total Mercury (may be combined with Total Metals)	EPA 7471	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	28 Days
Chromium, Hexavalent	EPA 7196	Aqueous	100mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> buffer	28 Days
Chromium, Hexavalent (subcontract)	EPA 7196	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	30 Days
Cyanide, Total	EPA 9014	Aqueous	125mL	125mL Polyethylene Bottle	Cool to $\leq 6^{\circ}\text{C}$ 1:1 NaOH to pH >8	14 Days
Cyanide, Total	EPA 9014	Solid	15g	4oz Glass Jar w/Teflon liner	Cool to $\leq 6^{\circ}\text{C}$	14 Days

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**Case Narrative**

**Lab # 53042 & 53078**

**Sample Receiving and Chain of Custody Discrepancies**

---

Samples were received in acceptable condition, between 0 and 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

**Calibration**

---

No exceptions noted.

**Method Blank**

---

No exceptions noted.

**Surrogate Recoveries**

---

Not applicable.

**Laboratory Control Sample Results**

---

No exceptions noted.

**Matrix Spike/Matrix Spike Duplicate/Duplicate Results**

---

Not requested for this project.

**Other**

---

No other exceptions noted.

**MassDEP Analytical Protocol Certification Form Questions A through I**

---

No explanation is needed for Questions A through I answered in the affirmative.

**Question G:** The CAM protocol reporting limits were not achieved for this project due to dilutions necessary for sample analysis. Box G is "No."

**Question H:** A solid MS/D was not specified. Box H is "No."

**Question I:** Metals: The MCP required metals were not requested by the customer. Box I is "No."

# GLOSSARY

%R	Percent Recovery
BLK	Blank (Method Blank, Preparation Blank)
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRM	Certified Reference Material (associated with solid Metals samples)
CRMD	Certified Reference Material Duplicate (associated with solid Metals samples)
Dil'n	Dilution
DL	Detection Limit
DUP	Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
LOD	Limit of Detection
LOQ	Limit of Quantitation
MB	Methanol Blank (associated with solid VOC samples)
MLCS	Methanol Laboratory Control Sample (associated with solid VOC samples)
MLCSD	Methanol Laboratory Control Sample Duplicate (associated with solid VOC samples)
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PB	Preparation Blank
QC	Quality Control
RL	Reporting Limit
RPD	Relative Percent Difference
SUR	Surrogate



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- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6020A	BLK12797	Lead		< 2.5	ug/g					
SW3051A6020A	CRM12797	Lead		175	ug/g	248		152 273		
SW3051A6020A	CRMD12797	Lead		180	ug/g	248		152 273	3	20
SW3051A6020A	MS12797	Lead	53062-009	380	ug/g	235	107	75 125		
SW3051A6020A	MSD12797	Lead	53062-009	350	ug/g	227	98	75 125	8	20





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 603-436-2001

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**CHAIN-OF-CUSTODY RECORD  
 AND ANALYSIS REQUEST**

**52513**

**ANALYSIS REQUEST**

Company Name: \_\_\_\_\_  
 Company Address: **SEE PAGE 1**  
 Report To: \_\_\_\_\_  
 Phone #: \_\_\_\_\_  
 Invoice to: \_\_\_\_\_  
 Email: **SEE PAGE 1**  
 PO #: \_\_\_\_\_

Project Name: \_\_\_\_\_  
 Project #: **SEE PAGE 1**  
 Project Location: NH MA ME VT \_\_\_\_\_  
 Accreditation Required? N/Y: \_\_\_\_\_  
 Protocol: RCRA SDWA NPDES  
           MCP NHDES DOD  
 Reporting QAPP GW-1 S-1  
 Limits: EPA DW Other  
 Quote # **N/A**  
 NH Reimbursement Pricing

<input checked="" type="checkbox"/> VOC 8260	<input type="checkbox"/> VOC 8260 MADEP *	<input type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MADEP *
<input type="checkbox"/> VOC 624.1	<input type="checkbox"/> VOC BTEX MBE, only	<input type="checkbox"/> VOC 8021VT	<input type="checkbox"/> VOC 8021VT
<input checked="" type="checkbox"/> TPH MADEP	<input type="checkbox"/> GPD 8015	<input type="checkbox"/> 1,4-Dioxane *	<input type="checkbox"/> 1,4-Dioxane *
<input type="checkbox"/> VOC 524.2	<input type="checkbox"/> VOC 524.2 NH List	<input type="checkbox"/> Gases-List	<input type="checkbox"/> Gases-List
<input type="checkbox"/> TPH	<input type="checkbox"/> DRD 8015	<input checked="" type="checkbox"/> EPH MADEP	<input type="checkbox"/> TPH Fingerprint
<input type="checkbox"/> 8270PAH	<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625.1	<input type="checkbox"/> EDB
<input type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608.3 Pest/PCB	<input type="checkbox"/> 608.3 Pest/PCB
<input type="checkbox"/> OAG 1664	<input type="checkbox"/> Mineral D&G 1664	<input type="checkbox"/> pH	<input type="checkbox"/> BOD
<input type="checkbox"/> pH	<input type="checkbox"/> Conductivity	<input type="checkbox"/> Turbidity	<input type="checkbox"/> Apparent Color
<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TS	<input type="checkbox"/> TVS
<input type="checkbox"/> RCRA Metals	<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input type="checkbox"/> Hardness
<input checked="" type="checkbox"/> Heavy Metals-list: <b>MCP 14 METALS</b>	<input type="checkbox"/> Dissolved Metals-list:	<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD
<input type="checkbox"/> Ammonia	<input type="checkbox"/> TOC	<input type="checkbox"/> TOX	<input type="checkbox"/> TN
<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Bacteria P/A	<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> Enterococci
<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide	<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Ormo P
<input type="checkbox"/> Nitrate	<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Sulfate
<input type="checkbox"/> Corrosivity	<input type="checkbox"/> Ignitability/FP	<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC
<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC	<input type="checkbox"/> TCLP Pesticide
<input type="checkbox"/> Subcontract	<input type="checkbox"/> Grain Size	<input type="checkbox"/> Herbicides	<input type="checkbox"/> Asbestos
<input type="checkbox"/> PFAS	<input type="checkbox"/> PFAS	<input type="checkbox"/> PFAS	<input type="checkbox"/> PFAS

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
525-13-12	SB-110(0-1')		X								4/7/20	12:45	AD
13	SB-110(0-3')		X								12:55	12:55	AD
14	SB-110(5-8')		X								12:55	12:50	AD
15	SB-112(0-1')		X								13:15		
16	SB-112(0-3')		X								13:20		
17	SB-112(5-8')		X								13:25		
18	MS-1		X								10:15		
19	MSD-1		X						X				
20	DUP-1		X										
21	TRIP BLANK												

**TAT REQUESTED**  
 Priority (24 hr)   
 Expedited (48 hr)   
 Standard (10 Business Days)  **5 DAY**  
 \*Date Needed

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
 \*VOCs & VPIT EXCEPT FOR TRIP BLANK  
 \*\*VPIT WITH TARGET VOCs

REPORTING INSTRUCTIONS  PDF (e-mail address)  EDD  
 HARD COPY REQUIRED  EDD

RECEIVED ON ICE  YES  NO  
 TEMPERATURE **2** °C

<b>CUSTODY RECORD</b> OSD-01 Revision 11/06/19	Relinquished by Sampler: _____	Date: 4/7/20	Time: 1500	Received by: _____	Date: 4-7	Time: 3:00
	Relinquished by: _____	Date: 4-7	Time: 16:00	Received by: _____	Date:	Time:
	Relinquished by: _____	Date:	Time:	Received by Laboratory: _____	Date:	Time:



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 603-436-2001  
 absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
 AND ANALYSIS REQUEST**

**53078**

**ANALYSIS REQUEST**

Company Name: **WESTON & SAMPSON**  
 Company Address: **55 WALKERS BROOK DR READING, MA**  
 Report To: **SARAH DESTEFANO & JILL MURPHY**  
 Phone #: **1 800 SAMPSON**  
 Invoice to: **SARAH DESTEFANO**  
 Email: **DESTEFANOS@WSEINC.COM**  
 PO #: **FREMINGTON BROWNFIELDS**

Project Name: **CEDAR WOODS**  
 Project #: **2180311**  
 Project Location: NH (MA) ME VT  
 Accreditation Required? N/Y: **Y**  
 Protocol: **RCRA SDWA NPDES**  
**MCP NHDES DOD**  
 Reporting: **SAPP 1 SW-1 (S-1)**  
 Limits: **EPA DW Other**  
 Quote #: **N/A**  
 NH Reimbursement Pricing

<input type="checkbox"/> VOC 8260	<input type="checkbox"/> VOC 8260 NHDES	<input type="checkbox"/> VOC 8260 MADEP	<input type="checkbox"/> VOC 8260
<input type="checkbox"/> VOC 624.1	<input type="checkbox"/> VOC BTEX MIBE only	<input type="checkbox"/> VOC 8021VT	<input type="checkbox"/> VOC 624.1
<input checked="" type="checkbox"/> VPH MADEP	<input type="checkbox"/> GRO 8015	<input checked="" type="checkbox"/> 1,4-Dioxane	<input checked="" type="checkbox"/> VPH MADEP
<input type="checkbox"/> VOC 524.2	<input type="checkbox"/> VOC 524.2 RH Lut	<input type="checkbox"/> Gases-Lut	<input type="checkbox"/> VOC 524.2
<input type="checkbox"/> TPH	<input checked="" type="checkbox"/> DRG 8015	<input checked="" type="checkbox"/> EPH MADEP	<input type="checkbox"/> TPH
<input type="checkbox"/> 8270PAH	<input type="checkbox"/> 8270ABN	<input type="checkbox"/> 625.1	<input type="checkbox"/> 8270PAH
<input type="checkbox"/> 8082 PCB	<input type="checkbox"/> 8081 Pesticides	<input type="checkbox"/> 608.3 Pest/PCB	<input type="checkbox"/> 8082 PCB
<input type="checkbox"/> O&G 1664	<input type="checkbox"/> Mineral O&G 1664		<input type="checkbox"/> O&G 1664
<input type="checkbox"/> PH	<input type="checkbox"/> BOD	<input type="checkbox"/> Conductivity	<input type="checkbox"/> PH
<input type="checkbox"/> TSS	<input type="checkbox"/> TDS	<input type="checkbox"/> TVS	<input type="checkbox"/> TSS
<input type="checkbox"/> RCRA Metals	<input type="checkbox"/> Priority Pollutant Metals	<input type="checkbox"/> TAL Metals	<input type="checkbox"/> RCRA Metals
<input checked="" type="checkbox"/> Total Metals-list: <b>MCP 14 METALS</b>			<input checked="" type="checkbox"/> Total Metals-list: <b>MCP 14 METALS</b>
<input type="checkbox"/> Dissolved Metals-list:			<input type="checkbox"/> Dissolved Metals-list:
<input type="checkbox"/> Ammonia	<input type="checkbox"/> COD	<input type="checkbox"/> TN	<input type="checkbox"/> Ammonia
<input type="checkbox"/> T-Phosphorus	<input type="checkbox"/> Bacteria P/A	<input type="checkbox"/> Bacteria MPN	<input type="checkbox"/> T-Phosphorus
<input type="checkbox"/> Cyanide	<input type="checkbox"/> Sulfide	<input type="checkbox"/> Nitrate + Nitrite	<input type="checkbox"/> Cyanide
<input type="checkbox"/> Nitrate	<input type="checkbox"/> Nitrite	<input type="checkbox"/> Chloride	<input type="checkbox"/> Nitrate
<input type="checkbox"/> Corrosivity	<input type="checkbox"/> Ignitibility/FP		<input type="checkbox"/> Corrosivity
<input type="checkbox"/> TCLP Metals	<input type="checkbox"/> TCLP VOC	<input type="checkbox"/> TCLP SVOC	<input type="checkbox"/> TCLP Metals
<input type="checkbox"/> Subcontract	<input type="checkbox"/> Grain Size	<input type="checkbox"/> Herbicides	<input type="checkbox"/> Subcontract
			<b>HEXAVALENT CHROMIUM (CR VI)</b>
			<b>Total Pb per Spacer 5741</b>

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix			Preservation Method				Sampling		SAMPLER
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	
52419-01	SB-106(0-1)		X			X				4/6/20	0920	
-02	SB-106(0-3)		X			X					0940	
-03	SB-106(5-8)		X			X					1000	
-04	SB-106(3-4)		X			X					0930	
-05	SB-105(0-3)		X			X					1040	
-06	SB-105(5-8)		X			X					1100	
-07	SB-102(0-3)		X			X					1145	
-08	SB-102(0-1)		X			X					1130	
-09	SB-102(11-14)		X			X					1140	
-10	SB-101(0-3)		X			X					1255	
-11	SB-101(9-12)		X			X					1305	

**TAT REQUESTED**  
 Priority (24 hr)   
 Expedited (48 hr)   
 Standard (10 Business Days)   
 \*Date Needed **5-DAY**

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

**SPECIAL INSTRUCTIONS**  
 \*VPIT NI TARGET VOCs

**REPORTING INSTRUCTIONS**  
 PDF (e-mail address) **murphy.j@wseinc.com**

HARD COPY REQUIRED  EDD

RECEIVED ON ICE  YES  NO  
 TEMPERATURE **0** °C

**CUSTODY RECORD**  
 OSD-01 Revision 11/06/19

Relinquished by:	Date: <b>4/6/20</b> Time: <b>12:40</b>	Received by:	Date: <b>4-6</b> Time: <b>1:41</b>
Relinquished by:	Date: <b>4-6</b> Time: <b>15:43</b>	Received by:	Date:   Time
Relinquished by:	Date:   Time	Received by Laboratory:	Date: <b>4/6/20</b> Time: <b>15:43</b>

Relayed Per J. Spencer 5/20/20

**Absolute Resource**  
associates



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Portsmouth, NH 03801  
603-436-2001

absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD  
AND ANALYSIS REQUEST**

**52498**

**ANALYSIS REQUEST**

Company Name: \_\_\_\_\_  
 Company Address: **SEE PAGE 1 OF 2**  
 Report To: \_\_\_\_\_  
 Phone #: \_\_\_\_\_  
 Invoice to: \_\_\_\_\_  
 Email: \_\_\_\_\_  
 PO #: \_\_\_\_\_

Project Name: \_\_\_\_\_  
 Project #: \_\_\_\_\_  
 Project Location: NH MA ME VT \_\_\_\_\_  
 Accreditation Required? N/Y: \_\_\_\_\_  
 Protocol: RCRA SDWA NPDES  
 MCP NHDES DOD  
 Reporting QAPP GW-1 S-1  
 Limits: EPA DW Other \_\_\_\_\_  
 Quote # \_\_\_\_\_  
 NH Reimbursement Pricing

VOC 8260  VOC 8260 NHDES  VOC 8260 MADEP  
 VOC 624.1  VOC BTEX MBE only  VOC 8021VT  
 ACPH MADEP  GRO 8015  1,4-Dioxane \*  
 VOC 524.2  VOC 524.2 NH List  Gases-List  
 TPH  DRO 8015  XEPH MADEP  TPH Fingerprint  
 8270PAH  8270A9H  625.1  EDB  
 8082 PCB  8081 Pesticides  608.3 Pest/PCB  
 O&G 1664  Mineral O&G 1664  
 pH  BOD  Conductivity  Turbidity  Apparent Color  
 TSS  TS  TVS  Alkalinity  Acidity  
 RCRA Metals  Priority Pollutant Metals  TAL Metals  Hardness  
 Total Metals-list: **MCP 14 METALS**  
 Dissolved Metals-list  
 Ammonia  COD  TKN  TN  TON  TOC  Ferrous Iron  
 T-Phosphorus  Bacteria P/A  Bacteria MPN  Enterococci  
 Cyanide  Sulfide  Nitrate + Nitrite  Ortho P  Phenols  
 Nitrate  Nitrite  Chloride  Sulfate  Bromide  Fluoride  
 Corrosivity  Ignitability/FP  
 TCLP Metals  TCLP VOC  TCLP SVOC  TCLP Pesticide  
 Subcontract:  Grain Size  Herbicides  Asbestos  PFAS  
**XX HEXAVALENT CHROMIUM (CR VI)**

Lab Sample ID <small>(Lab Use Only)</small>	Field ID	# CONTAINERS	Matrix			Preservation Method					Sampling		
			WATER	SOLID	OTHER	HCl	HNO <sub>3</sub>	H <sub>2</sub> SO <sub>4</sub>	NaOH	MeOH	DATE	TIME	SAMPLER
52498-12	SB-103(0-3)			XX							4/6/20	1220	
-13	SB-103(3-5)			XX							↓	1225	
-14	SB-103(7-10)			XX								1225	
-15	TREP BLANK			X							4/6/20		X

**TAT REQUESTED**  
 Priority (24 hr)   
 Expedited (48 hr)   
 Standard (10 Business Days)   
 \*Date Needed **5-DAY**

**SPECIAL INSTRUCTIONS**  
 \*VPH WITH TARGET VOCs

**CUSTODY RECORD**  
QSD-01 Revision 11/06/19

Relinquished by: \_\_\_\_\_  
 Date: 4/6/20 Time: 1340  
 Relinquished by: \_\_\_\_\_  
 Date: 4-6 Time: 1543  
 Relinquished by: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_

Received by: \_\_\_\_\_  
 Date: 4-6 Time: 1-40  
 Received by: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 Received by Laboratory: \_\_\_\_\_  
 Date: 4/6/20 Time: 1543

RECEIVED ON ICE  YES  NO  
 TEMPERATURE 0 °C

MDM 4/6/20  
 GG Jc



APPENDIX C

Risk Characterization Tables

## Method 3 Imminent Hazard Assessment for Park Visitor Exposed to Chemicals in Soil - Shortform 2012 (sf12psih)

### Index

#### Tab

<b>EPCs</b>	Table PSIH-1: Select chemicals and enter Exposure Point Concentrations (EPCs). Estimated risks are presented to the right.
<b>C Eq</b>	Table PSIH-2: Equations to calculate cancer risks.
<b>NC Eq</b>	Table PSIH-3: Equations to calculate noncancer risks.
<b>Exp</b>	Table PSIH-4: Definitions and exposure factors.
<b>Chem</b>	Table PSIH-5: Chemical-specific data.
<b>Cyanide</b>	Table PSIH-6: Cyanide Calculation

---

Spreadsheets designed by Rafael McDonald, MassDEP

Questions and Comments may be addressed to:

**Lydia Thompson**

Massachusetts Department of Environmental Protection

Office of Research and Standards

One Winter Street

Boston, MA 02108 USA

Telephone: (617) 556-1165

Fax: (617) 556-1006

Email: [Lydia.Thompson@state.ma.us](mailto:Lydia.Thompson@state.ma.us)

---

**Park Visitor - Soil Imminent Hazard Evaluation: Table PSIH-1**  
**Exposure Point Concentration (EPC)**  
**Based on Visitor Ages 1-6 (Cancer) and 1-2 (Noncancer)**

ShortForm Version 10-12

Vlookup Version v0315

**\*\*Do not insert or delete any rows\*\***

Click on empty cell below and select OHM using arrow.

ELCR (all chemicals) =  
 Subchronic HI (all chemicals) = 2.4E+00

Oil or Hazardous Material	EPC (mg/kg)	ELCR <sub>ingestion</sub>	ELCR <sub>dermal</sub>	ELCR <sub>total</sub>	Subchronic		
					HQ <sub>ing</sub>	HQ <sub>derm</sub>	HQ <sub>total</sub>
LEAD	8.5E+02				2.3E+00	1.6E-01	2.4E+00

**Note! Lead IH HQ limit is 1, not 10.**

**Park Visitor - Soil: Table PSIH-2**  
**Equations to Calculate Cancer Risk for Visitor (Age 1-6 years)**

Vlookup Version v0315

**Cancer Risk from Ingestion**

$$ELCR_{ing} = LADD_{ing} * CSF$$

$$LADD_{ing} = \frac{[OHM]_{soil} * IR_x * RAF_{c-ing} * EF_{ing} * ED * EP * C}{BW * AP_{lifetime}}$$

**Cancer Risk from Dermal Absorption**

$$ELCR_{derm} = LADD_{derm} * CSF$$

$$LADD_{derm} = \frac{[OHM]_{soil} * SA * RAF_{c-derm} * SAF * EF_{derm} * ED * EP * C}{BW * AP_{lifetime}}$$

Parameter	Value	Units
CSF	OHM specific	(mg/kg-day) <sup>-1</sup>
LADD	age/OHM specific	mg/kg-day
[OHM] <sub>soil</sub>	OHM specific	mg/kg
IR	100	mg/day
RAF <sub>c-ing</sub>	OHM specific	dimensionless
RAF <sub>c-derm</sub>	OHM specific	dimensionless
EF <sub>ing,derm</sub>	0.247	event/day
ED	1	day/event
EP	5	years
C	0.000001	kg/mg
BW	14.6	kg
AP <sub>(lifetime)</sub>	70	years
SA	2231	cm <sup>2</sup> /day
SAF	0.35	mg/cm <sup>2</sup>

**Park Visitor - Soil: Table PSIH-3**  
**Equations to Calculate Noncancer Risk for Visitor (Age 1-2 years)**

Vlookup Version v0315

**Noncancer Risk from Ingestion**

$$HQ_{ing} = \frac{ADD_{ing}}{RfD_{subchronic}}$$

$$ADD_{ing} = \frac{[OHM]_{soil} * IR * RAF_{nc-ing} * EF_{ing} * ED * EP * C}{BW * AP}$$

**Noncancer Risk from Dermal Absorption**

$$HQ_{derm} = \frac{ADD_{derm}}{RfD_{subchronic}}$$

$$ADD_{derm} = \frac{[OHM]_{soil} * SA * RAF_{nc-derm} * SAF * EF_{derm} * ED * EP * C}{BW * AP}$$

Parameter	Value	Units
RfD	OHM specific	mg/kg-day
ADD	OHM specific	mg/kg-day
[OHM] <sub>soil</sub>	OHM specific	mg/kg
IR	100	mg/day
RAF <sub>nc-ing</sub>	OHM specific	dimensionless
RAF <sub>nc-derm</sub>	OHM specific	dimensionless
EF <sub>ing,derm</sub>	0.428	event/day
ED	1	day/event
EP	0.577	years
C	0.000001	kg/mg
BW	10.7	kg
AP	0.577	year
SA	1670	cm <sup>2</sup> /day
SAF	0.35	mg/cm <sup>2</sup>

**Park Visitor - Soil: Table PSIH-4  
Definitions and Exposure Factors**

Vlookup Version v0315

Parameter	Value	Units	Notes
ELCR - Excess Lifetime Cancer Risk	chemical specific	dimensionless	Pathway specific (ing =ingestion, derm=dermal)
CSF - Cancer Slope Factor	chemical specific	(mg/kg-day) <sup>-1</sup>	see Table PSIH-5.
LADD - Lifetime Average Daily Dose	chemical specific	mg/kg-day	Pathway specific - see Table PSIH-2.
HQ - Hazard Quotient	chemical specific	dimensionless	Pathway specific (ing =ingestion, derm=dermal) - see Table PSIH-3.
RfD - Reference Dose	chemical specific	mg/kg-day	see Table PSIH-5.
ADD - Average Daily Dose	chemical specific	mg/kg-day	Pathway specific
EPC - Exposure Point Concentration	chemical specific	mg/kg	
IR - Soil Ingestion Rate	100	mg/day	
RAF <sub>c</sub> - Relative Absorption Factor for Cancer Effects	chemical specific	dimensionless	MADEP. 1995. Guidance for Disposal Site Risk Characterization. Appendix Table B-3.
RAF <sub>NC</sub> - Relative Absorption Factor for non-Cancer Effects	chemical specific	dimensionless	Adjusts estimated dose to conform to the relevant CSF. See Table PS-6
EF <sub>subchronic</sub> - Exposure Frequency for subchronic exposure	0.428	event/day	3 events/week
EF <sub>lifetime</sub> - Exposure Frequency for chronic or lifetime exposure	0.247	event/day	3 events/week, 30 weeks/year
ED - Exposure Duration	1	day/event	
EP <sub>(1-2)</sub> - Exposure Period for age group 1-2	0.577	years	30 weeks
EP <sub>(1-6)</sub> - Exposure Period for age group 1-6	5	years	
BW <sub>(1-2)</sub> - Body Weight for age group 1-2	10.7	kg	U.S. EPA. 1997. Exposure Factors Handbook. Table 7-7, females.
BW <sub>(1-6)</sub> - Body Weight for age group 1-6	14.6	kg	Ibid
AP <sub>subchronic</sub> - Averaging Period for subchronic noncancer	0.577	years	30 weeks
AP <sub>lifetime</sub> - Averaging Period for cancer/lifetime	70	years	
SA <sub>(1-2)</sub> - Surface Area for age group 1-2	1670	cm <sup>2</sup> /day	50th percentile of face (1/3 head), forearms, hands, lower legs, and feet for females MADEP 1995 Guidance for Disposal Site Risk Characterization, Appendix Table B-2.
SA <sub>(1-6)</sub> - Surface Area for age group 1-6	2231	cm <sup>2</sup> / day	Ibid
SAF - Surface Adherence Factor	0.35	mg <sub>soil</sub> / cm <sup>2</sup>	All SAFs developed for ShortForm according to procedure outlined in MA DEP Technical Update: Weighted Skin-Soil Adherence Factors, April 2002

**Park Visitor - Soil: Table PSIH-5  
Chemical-Specific Data**

Vlookup Version v0315

Oil or Hazardous Material	CSF (mg/kg-day) <sup>-1</sup>	RAF <sub>c-ing</sub>	RAF <sub>c-derm</sub>	Subchronic RfD mg/kg-day	Subchronic RAF <sub>nc-ing</sub>	Subchronic RAF <sub>nc-derm</sub>
LEAD				7.5E-04	0.5	0.006

APPENDIX D  
Photographic Log



Immediate Response Action (IRA) Plan  
618R Waverly Street, Framingham, MA  
Release Tracking Number (RTN) 3-36304

PHOTO DOCUMENTATION



Photo 1

View looking south at the northeastern corner of the temporary fence installed at the Site.



Photo 2

View looking south at the northern portion of the Site, from the paved parking lot to the north.



Photo 3

View looking south at the northwestern corner of the temporary fence installed at the Site. The posted signage is visible in the center of the photo.

## APPENDIX E

### IDW Disposal Documentation



# DAILY WORKSHEET

3

Date: 7/29/20 Project Number: 20-0020

CLIENT / SITE INFORMATION Weston & Sampson

Name: Cedar Woods

Address: 618 R Waverly St, Framingham, MA

Contact: Joe Spencer 781-443-2779

PROJECTS NOTES

1-5011 bearings

Pick up Drum Dispose @ Northland

LABOR onsite 1145-1200

MATERIALS / EQUIPMENT / TOOLS

Name	Position	Travel	On-site	Off-Site	Travel	Quantity	Item / Description
<u>M. Lynch</u>	<u>ES</u>	<u>1.5</u>	<u>.25</u>		<u>1.25</u>	<u>1</u>	<u>PPE Level: A B C (D)</u>
						<u>1</u>	<u>Box truck</u>
						<u>1</u>	<u>waste label</u>

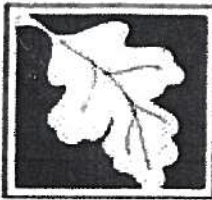
SUBCONTRACTORS

Disposal @ Northland

WEATHER OBSERVATIONS

Project Manager's Signature: \_\_\_\_\_

Client's Signature: \_\_\_\_\_



**BILL OF LADING (pursuant to 310 CMR 40.0030)**

**A. LOCATION OF SITE OR DISPOSAL SITE WHERE REMEDIATION WASTE WAS GENERATED:**

1. Release Name/Location Aid: PROPERTY TO EAST OF WAVERLY & MELLENS STS
2. Street Address: 618R WAVERLY STREET
3. City/Town: FRAMINGHAM 4. Zip Code: 017020000
5. Check here if the disposal site that is the source of the release is Tier Classified. Check the current Tier Classification Category.  
 a. Tier I     b. Tier ID     c. Tier II

**B. THIS FORM IS BEING USED TO: (check one: B1-B4):**

1. Submit a **Bill of Lading (BOL)** to transport Remediation Waste to Temporary Storage or a Receiving Facility.  
 Response Actions associated with this BOL (check all that apply):  
 a. Immediate Response Action (IRA)     e. Comprehensive Response Actions  
 b. Release Abatement Measure (RAM)     f. Limited Removal Action (LRA): (must be retained pursuant to 310 CMR 40.0034(6); can't be submitted via eDEP)  
 c. Downgradient Property Status (DPS)     g. Other \_\_\_\_\_  
 d. Utility Release Abatement Measure (URAM)
2. Submit an Attestation of Completion of Shipment to Temporary Storage (Sections C, F and J are not required):
3. Submit an Attestation of Completion of Shipment to a Receiving Facility (Sections C, F and J are not required):
4. Certify that Remediation Waste Was Not Shipped, and the Bill of Lading is Void. (Sections C, D, E, and F are not required)
5. Date Bill of Lading submitted to the Department: \_\_\_\_\_ b. eDEP Transaction ID: \_\_\_\_\_  
 (mm/dd/yyyy)
6. Period of Generation Associated with this Bill of Lading 4/6/2020 to 4/8/2020  
 (mm/dd/yyyy) (mm/dd/yyyy)

(All sections of this transmittal form must be filled out unless otherwise noted above)

The Bill of Lading is not considered complete until the Attestation of Completion of Shipment is received by the Department.

**C. DESCRIPTION OF WASTE AND WASTE SOURCE:**

1. Contaminated Media/Debris (check all that apply):  
 a. Soil     b. Groundwater     c. Surface Water     d. Sediment     e. Vegetation or Organic Debris  
 f. Demolition/Construction Waste     g. Inorganic Absorbent Materials     h. Other: \_\_\_\_\_
2. Uncontainerized Waste (check all that apply):  
 a. Inorganic Absorbent Materials     b. Other: \_\_\_\_\_



C. DESCRIPTION OF WASTE AND WASTE SOURCE (cont.):

3. Containerized Waste (check all that apply):

- a. Tank Bottoms/Sludges
- b. Containers
- c. Drums
- d. Engineered Impoundments
- e. Other: \_\_\_\_\_

4. Estimated Quantity: 0.2  Tons  Cu. Yds.  Gallons

5. Contaminant Source (check one):

- a. Transportation Accident
- b. Underground Storage Tank
- c. Brownfields Redevelopment
- d. Other: URBAN FILL/IMPACTED FILL MATERIAL

6. Type of Contaminant (check all that apply):

- a. Gasoline
- b. Diesel Fuel
- c. #2 Fuel Oil
- d. #4 Fuel Oil
- e. #6 Fuel Oil
- f. Jet Fuel
- g. Waste Oil
- h. Kerosene
- i. Chlorinated Solvents
- j. Urban Fill
- k. Other: \_\_\_\_\_

7. Constituents of Concern (check all that apply):

- a. As
- b. Cd
- c. Cr
- d. Pb
- e. Hg
- f. EPH/TPH
- g. VPH
- h. PCBs
- i. VOCs
- j. SVOCs
- k. Other: ZN, SB

8. If applicable, check the box for the Reportable Concentration Category of the site:

- a. RCS-1
- b. RCS-2
- c. RCGW-1
- d. RCGW-2

9. Remediation Waste Characterization Documentation (check at least one):

- a. Site History Information
- b. Sampling Analytical Methods and Procedures
- c. Laboratory Data
- d. Field Screening Data
- e. Characterization Documentation previously submitted to the Department

i. Date submitted: \_\_\_\_\_ ii. Type of Documentation: \_\_\_\_\_  
(mm/dd/yyyy)

D. TRANSPORTER OR COMMON CARRIER INFORMATION:

1. Transporter/Common Carrier Name: STRATEGIC ENVL SERVICES

2. Contact First Name: ALISA 3. Last Name: SEELEY

4. Street: 27 KREIGER LANE UNIT 10 5. Title: WASTE DISPOSAL COORDINATOR

6. City/Town: GLASTONBURY 7. State: CT 8. Zip Code: 060330000

9. Telephone: 8602662616 10. Ext: \_\_\_\_\_ 11. Email: aseeley@strategic-es.com





**G. PERSON SUBMITTING BILL OF LADING:**

1. Check all that apply:  a. change in contact name  b. change of address  c. change in the person undertaking response actions
2. Name of Organization: CITY OF FRAMINGHAM
3. Contact First Name: THATCHER 4. Last Name: KEZER III
5. Street: 150 CONCORD STREET 6. Title: CHIEF OPERATING OFFICER
7. City/Town: FRAMINGHAM 8. State: MA 9. Zip Code: 017020000
10. Telephone: 5085325400 11. Ext: \_\_\_\_\_ 12. Email: tkezer@framinghamma.gov

**H. RELATIONSHIP TO SITE OF PERSON SUBMITTING BILL OF LADING:**

Check here to change relationship

1. RP or PRP  a. Owner  b. Operator  c. Generator  d. Transporter
- e. Other RP or PRP Specify: \_\_\_\_\_

2. Fiduciary, Secured Lender or Municipality with Exempt Status (as defined by M.G.L. c. 21E, s. 2)

3. Agency or Public Utility on a Right of Way (as defined by M.G.L. c. 21E, s. 5(j))

4. Any Other Person Undertaking Response Actions: Specify Relationship: \_\_\_\_\_

**I. REQUIRED ATTACHMENT AND SUBMITTALS:**

1. Check here if the Response Action(s) on which this opinion is based, if any, are (were) subject to any order(s), permit(s) and/or approvals issued by DEP or EPA. If the box is checked, you must attach a statement identifying the applicable provisions thereof.
2. Check here if any non-updatable information provided on this form is incorrect, e.g. Release Address/Location Aid. Send corrections to BWSC.eDEP@state.ma.us
3. Check here to certify that the LSP Opinion containing the material facts, data, and other information is attached.

**J. CERTIFICATION OF PERSON SUBMITTING BILL OF LADING:**

1. I, THATCHER W. KEZER III, attest under the pains and penalties or perjury (i) that I have personally examined and am familiar with the information contained in this submittal, including any and all documents accompanying this transmittal form, (ii) that, based on my inquiry of those individuals immediately responsible for obtaining the information, the material information contained in this submittal is, to the best of my knowledge and belief, true, accurate and complete, and (iii) that I am fully authorized to make this attestation on behalf of the entity legally responsible for this submittal. I/the person or entity on whose behalf this submittal is made am/is aware that there are significant penalties, including, but not limited to, possible fines and imprisonment, for willfully submitting false, inaccurate, or incomplete information.

2. By: THATCHER W. KEZER III 3. Title: CHIEF OPERATING OFFICER
4. For: CITY OF FRAMINGHAM 5. Date: 7/28/2020  
(Name of person or entity recorded in Section G) (mm/dd/yyyy)



**J. CERTIFICATION OF PERSON SUBMITTING BILL OF LADING (cont.) :**

6. Check here if the address of the person providing certification is different from address recorded in Section G.

7. Street: \_\_\_\_\_

8. City/Town: \_\_\_\_\_ 9. State: \_\_\_\_\_ 10. Zip Code: \_\_\_\_\_

11. Telephone: \_\_\_\_\_ 12. Ext: \_\_\_\_\_ 13. Email: \_\_\_\_\_

**YOU ARE SUBJECT TO AN ANNUAL COMPLIANCE ASSURANCE FEE OF UP TO \$10,000 PER BILLABLE YEAR FOR THIS DISPOSAL SITE. YOU MUST LEGIBLY COMPLETE ALL RELEVANT SECTIONS OF THIS FORM OR DEP MAY RETURN THE DOCUMENT AS INCOMPLETE. IF YOU SUBMIT AN INCOMPLETE FORM, YOU MAY BE PENALIZED FOR MISSING A REQUIRED DEADLINE.**

Date Stamp (MassDEP USE ONLY):

Received by DEP on 7/28/2020 6:19:08 PM





**Massachusetts Department of Environmental Protection**  
**Bureau of Waste Site Cleanup**

**BILL OF LADING Transport Log Sheet**

Release Tracking Number

Page 1 OF 1

-

<b>I. LOAD INFORMATION:</b>	
<b>Load 1:</b> Signature of Transporter Representative: <i>[Signature]</i> Date of Shipment: <u>7/29/20</u> Time of Shipment: <u>1145</u> <input checked="" type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: <u>MA T33-466</u> Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: <i>[Signature]</i> Date of Receipt: <u>7/29/20</u> Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):
<b>Load 2:</b> Signature of Transporter Representative: Date of Shipment: Time of Shipment: <input type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: Date of Receipt: Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):
<b>Load 3:</b> Signature of Transporter Representative: Date of Shipment: Time of Shipment: <input type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: Date of Receipt: Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):
<b>Load 4:</b> Signature of Transporter Representative: Date of Shipment: Time of Shipment: <input type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: Date of Receipt: Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):
<b>Load 5:</b> Signature of Transporter Representative: Date of Shipment: Time of Shipment: <input type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: Date of Receipt: Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):
<b>Load 6:</b> Signature of Transporter Representative: Date of Shipment: Time of Shipment: <input type="checkbox"/> AM <input type="checkbox"/> PM Truck/Tractor Registration: Trailer Registration (if any):	Receiving Facility/Temporary Storage Representative: Date of Receipt: Time of Receipt: <input type="checkbox"/> AM <input type="checkbox"/> PM Load Size (cu. yds./tons):

**J. LOG SHEET VOLUME INFORMATION:**

Total Volume Recorded This Page (cu. yds./tons)
Total Carried Forward (cu. yds./tons):
Total Carried Forward and This Page (cu. yds./tons):

July 24, 2020

Massachusetts Department of Environmental Protection  
Northeast Regional Office  
Bureau of Waste Site Cleanup  
205B Lowell Street  
Wilmington, Massachusetts 01887

Re: **LSP Opinion - Soil Disposal**  
618R Waverly Street  
Framingham, Massachusetts  
RTN 3-36304

Dear Sir/Madam:

Weston & Sampson has prepared this Licensed Site Professional (LSP) Opinion letter for the disposal of one 55-gallon drum of soil that was generated during assessment activities at the above-referenced property, located in Framingham, Massachusetts (the Site). See Figure 1 for the general location of the Site.

The Site is regulated under the Massachusetts Contingency Plan (MCP), and the subject soil was generated during Site assessment activities conducted under the City of Framingham's US EPA Brownfields Grant. Based on the data obtained during Site assessment activities, a 2-hour Notification Condition was identified due to elevated concentrations of lead identified in a small area (approximately 500 sf) of shallow soil. On June 3, 2020, the City reported the Release to MassDEP, and MassDEP assigned Release Tracking Number (RTN) 3-36304 to the Release.

### **Current and Former Site Use**

The Site consists of approximately 2.08 acres of vacant land, comprised of woodlands, wetland, and paved parking lot. The 618R Waverly Street parcel (parcel ID: 134-64-7867) was initially part of a larger parcel (with current day 618 Waverly Street formerly known comprehensively as 612 Waverly Street), that was formerly occupied by Siltan Glass Co. Historical operations on the property included an auto parts retailer, filling station, and auto glass facility.

An ASTM Phase I ESA in 2009, ASTM Phase II ESA in 2010, and cleanup/reuse planning was conducted for the former Siltan Glass property (612 Waverly Street). An Activity and Use Limitation (AUL), for release tracking number (RTN) 3-29744, was filed in 2012 for the property by the then current property owner Charles L. Siltan, Inc. Contaminants of concern at the property included metals, Extractable Petroleum Hydrocarbons (EPH) and Polycyclic Aromatic Hydrocarbons (PAHs). These impacts were associated with historic fill materials and point source releases of petroleum from past operations as a gasoline station and salvage facility. The AUL restricts the use of the property to commercial / industrial uses only and requires a Soil Management Plan and Health and Safety Plan prior to any soil disturbance.

The Siltan Glass property was acquired in 2015 by the City of Framingham through tax title taking, and the property was subsequently subdivided to support sale of the existing facility structure as a commercial property (currently addressed as 618 Waverly St). Ownership of the remainder of the Siltan Glass property (currently addressed as 618R Waverly Street), was retained by the City of Framingham.

### **Soil Generation and Sampling**

From April 6 through 8, 2020, Weston & Sampson oversaw the advancement of 15 soil borings (SB-101 through SB-115) at locations depicted in Figure 2, Site Plan. During the course of drilling activities, approximately 40 gallons of surplus drill cuttings were placed in a labeled and secured 55-gallon steel drum and stored on-Site (see Figure 2 for location of drum). The soil was generated in small volumes from each of

the boring locations depicted on the attached Figure 2. Thirty-five soil samples were collected from discrete depth intervals, as shown on Tables 1A-1C and Table 2. Samples were submitted to Absolute Resource Associates of Portsmouth, NH (ARA) for analysis of antimony, arsenic, barium, beryllium, cadmium, chromium (III), chromium (VI), lead, mercury, nickel, silver, thallium, vanadium, zinc, extractable petroleum hydrocarbons (EPH) with target polycyclic aromatic hydrocarbons (PAHs), volatile petroleum hydrocarbons (VPH) with target volatile organic compounds (VOCs), and/or VOCs.

On June 15, 2020, Weston & Sampson collected one composite sample from the drummed soil (IDW-1) and submitted it to ARA for analysis of arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver by Toxicity Characteristic Leaching Procedures (TCLP) methods. Copies of the laboratory analytical reports are included as an attachment.

### **Analytical Results**

Data in the attached tables were compared to applicable MCP Reportable Concentrations criteria (RCS-1).

#### Metals

All metals, excepting selenium, were detected above laboratory reporting limits in at least one of the 35 soil samples submitted for analysis. Antimony, arsenic, lead, and zinc were detected at concentrations exceeding the applicable RCS-1 threshold in multiple samples. The result of the TCLP analysis for the subject sample showed concentrations below levels indicative of a characteristic hazardous waste.

#### EPH/PAHs

EPH aliphatic and aromatic fractions, as well as target PAHs were detected above the laboratory reporting limits at multiple boring locations. The PAHs benzo(a)pyrene and dibenz(a,h)anthracene were detected at concentrations exceeding the respective MCP RCS-1 in multiple samples.

#### VPH/VOCs

VPH aliphatic and aromatic fractions, as well as target petroleum VOCs were detected above the laboratory reporting limits, but below the respective RCS-1 threshold at three locations.

In addition, including two chlorinated VOCs (cis-1,2-dichloroethylene and trichloroethylene) were detected in sample SB-108 (6-9) at concentrations above the respective RCS-1 threshold.

### **Closure Statement**

Due diligence was used to characterize the soil for the presence of listed hazardous waste and characteristic hazardous waste in accordance with DEP Policy # HW93-01. Historical Site review and chemical testing performed on the soil did not indicate the presence of a listed hazardous waste nor did the soil exhibit a characteristic of hazardous waste. Importantly, there is no known history of activities involving cis-1,2-dichloroethylene and trichloroethylene that indicate the presence of these contaminants constitute a hazardous waste.

In consideration of the Site history and current conditions at the Site from which the soil proposed for disposal has been generated, it is our opinion that the samples collected to characterize the soil were sufficient in terms of sampling methods, sampling dates, analytes tested for, QA/QC criteria, number of samples, sampling depth, and sampling locations to provide analytical data which are representative of the soil proposed for such disposal/re-use.

Weston & Sampson has completed an evaluation of the laboratory analytical QA/QC data accompanying the laboratory reports and this submittal contains data which are judged to be useable and representative. Based on our review of the chemical soil testing data obtained during this sampling program, the soils represented by this submittal are suitable for disposal as non-hazardous waste at a landfill. If you have any questions or comments regarding this letter or need any additional information, please do not hesitate to contact our office at (978) 532-1900.

Sincerely,

WESTON & SAMPSON ENGINEERS, INC.

A handwritten signature in black ink, appearing to read "F. M. Ricciardi". The signature is fluid and cursive, with a large loop at the end.

Frank M. Ricciardi, PE, LSP  
Vice President

Attachments:

Figures

Tables

Stericycle Waste Profile Sheet

Laboratory Analytical Reports

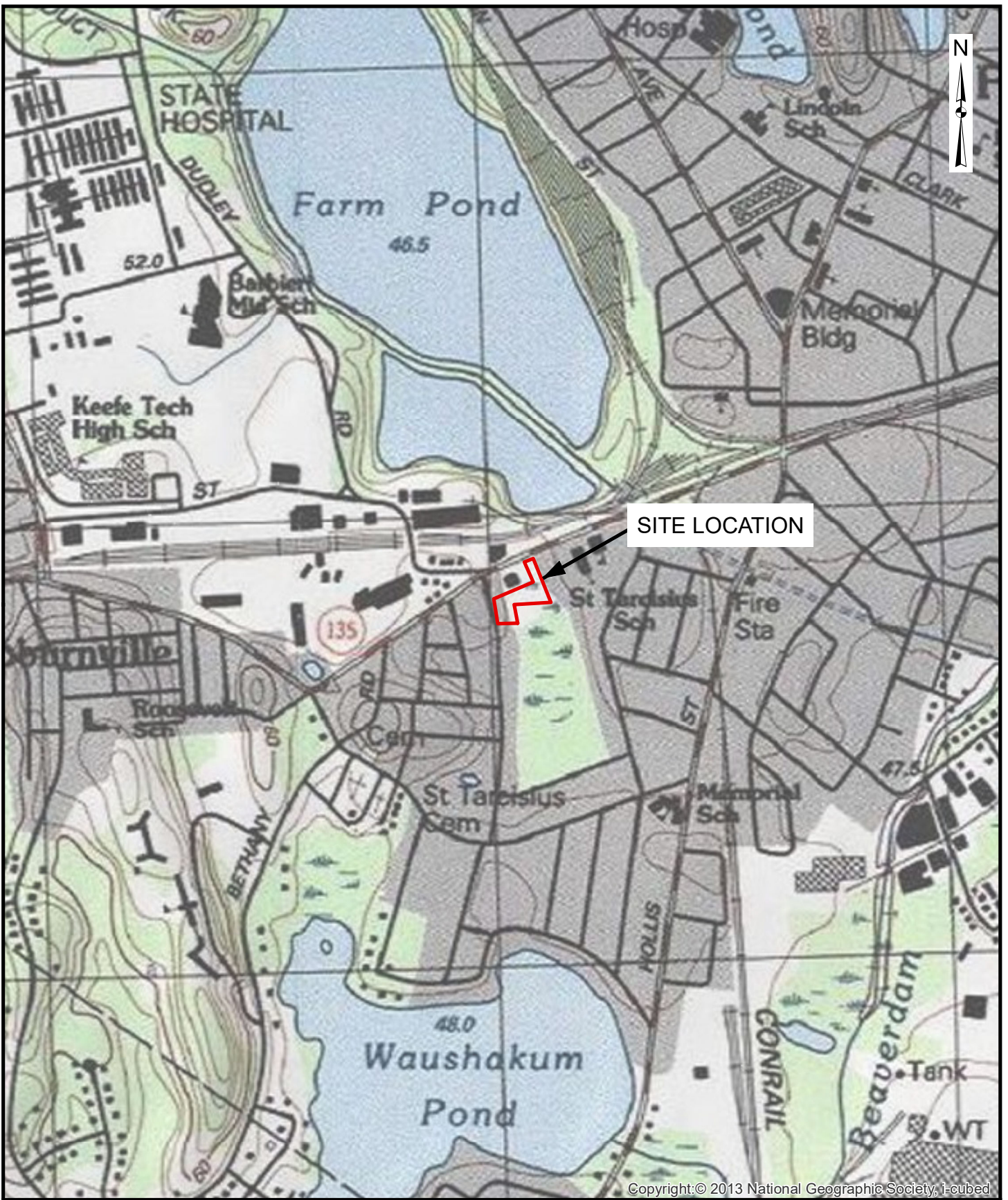
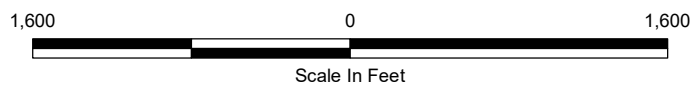
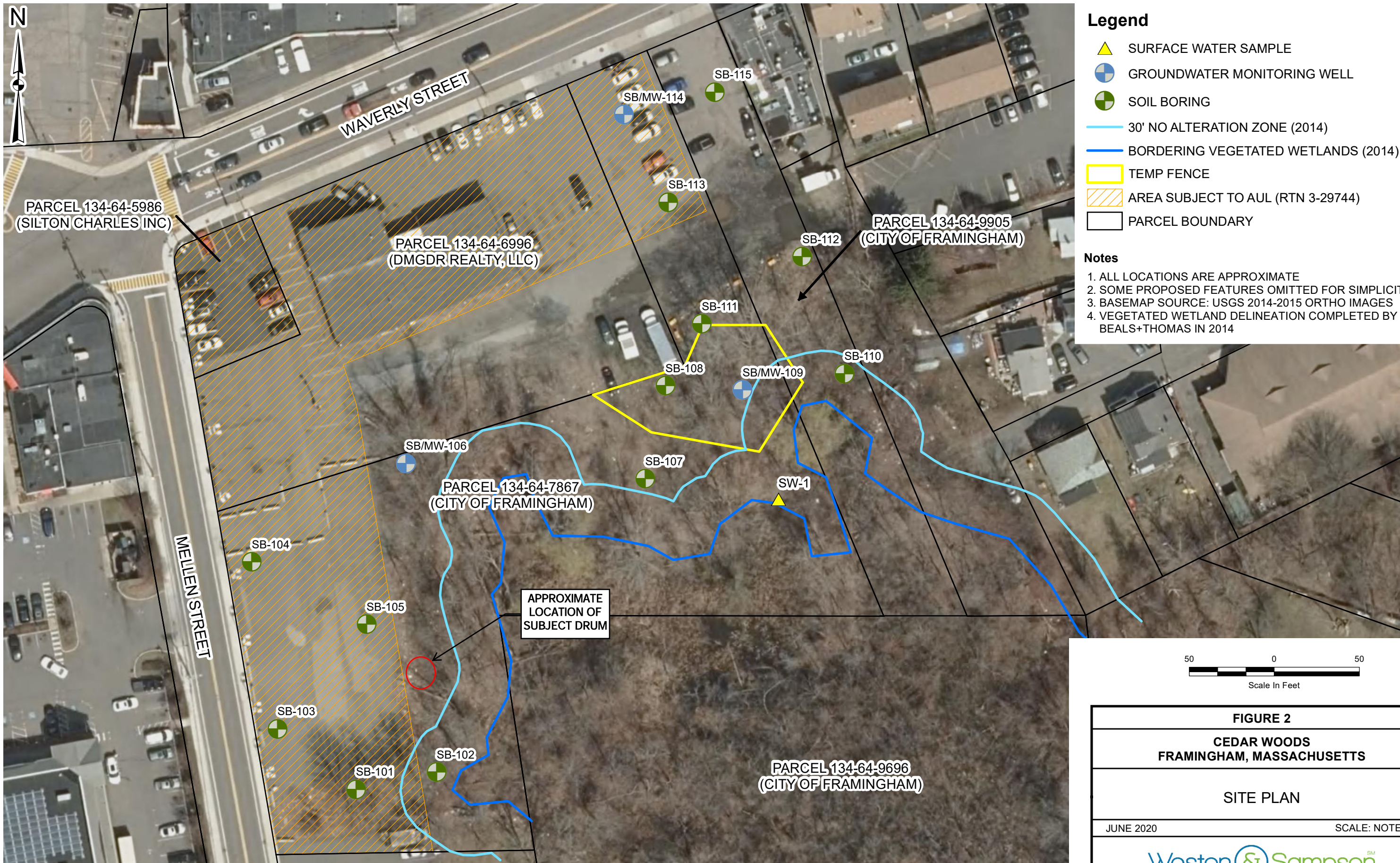


FIGURE 1  
 CEDAR WOODS  
 FRAMINGHAM, MA  
 LOCUS MAP



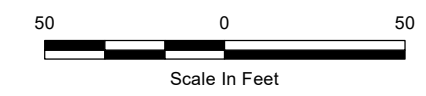


**Legend**

- ▲ SURFACE WATER SAMPLE
- ⊕ GROUNDWATER MONITORING WELL
- ⊕ SOIL BORING
- 30' NO ALTERATION ZONE (2014)
- BORDERING VEGETATED WETLANDS (2014)
- TEMP FENCE
- AREA SUBJECT TO AUL (RTN 3-29744)
- PARCEL BOUNDARY

**Notes**

1. ALL LOCATIONS ARE APPROXIMATE
2. SOME PROPOSED FEATURES OMITTED FOR SIMPLICITY
3. BASEMAP SOURCE: USGS 2014-2015 ORTHO IMAGES
4. VEGETATED WETLAND DELINEATION COMPLETED BY BEALS+THOMAS IN 2014



<b>FIGURE 2</b>	
<b>CEDAR WOODS FRAMINGHAM, MASSACHUSETTS</b>	
<b>SITE PLAN</b>	
JUNE 2020	SCALE: NOTED

**Table 1A**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**Cedar Woods**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)													
			RCS-1	S-1/GW-2	S-1/GW-3	SB-101		SB-102		SB-103			SB-104		SB-105	
						0-3 feet 4/6/2020	9-12 feet 4/6/2020	0-3 feet 4/6/2020	11-14 feet 4/6/2020	0-3 feet 4/6/2020	3-5 feet 4/6/2020	7-10 feet 4/6/2020	0-3 Feet 4/7/2020	8-10 Feet 4/7/2020	0-3 feet 4/6/2020	5-8 feet 4/6/2020
<b>EPH</b>																
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<21	<25	<20	<46	NT	NT	NT	<20	<22	<24	<22	
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>35</b>	<b>82</b>	<b>130</b>	<b>80</b>	NT	NT	NT	<b>34</b>	<b>100</b>	<b>280</b>	<b>23</b>	
C11-C22 Aromatics	mg/kg	1000	1000	1000	<21	<b>41</b>	<b>50</b>	<46	NT	NT	NT	<b>25</b>	<22	<b>130</b>	<22	
<b>Target PAHS</b>																
Acenaphthene	mg/kg	4	1000	1000	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Acenaphthylene	mg/kg	1	600	10	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Anthracene	mg/kg	1000	1000	1000	<0.21	<b>0.40</b>	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Benzo(A)Anthracene	mg/kg	7	7	7	<0.21	<b>0.98</b>	<0.20	<0.46	NT	NT	NT	<b>0.40</b>	<0.22	<b>1.5</b>	<0.22	
Benzo(A)Pyrene	mg/kg	2	2	2	<0.21	<b>0.89</b>	<0.20	<0.46	NT	NT	NT	<b>0.44</b>	<0.22	<b>1.4</b>	<0.22	
Benzo(B)Fluoranthene	mg/kg	7	7	7	<0.21	<b>0.88</b>	<0.20	<0.46	NT	NT	NT	<b>0.46</b>	<0.22	<b>1.6</b>	<0.22	
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<0.21	<b>0.72</b>	<0.20	<0.46	NT	NT	NT	<b>0.35</b>	<0.22	<b>1.2</b>	<0.22	
Benzo(K)Fluoranthene	mg/kg	70	70	70	<0.21	<b>0.71</b>	<0.20	<0.46	NT	NT	NT	<b>0.44</b>	<0.22	<b>1.2</b>	<0.22	
Chrysene	mg/kg	70	70	70	<0.21	<b>1.0</b>	<0.20	<0.46	NT	NT	NT	<b>0.56</b>	<0.22	<b>1.9</b>	<0.22	
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<b>0.38</b>	<0.22	
Fluoranthene	mg/kg	1000	1000	1000	<b>0.22</b>	<b>2.1</b>	<0.20	<b>0.48</b>	NT	NT	NT	<b>0.80</b>	<0.22	<b>3.2</b>	<0.22	
Fluorene	mg/kg	1000	1000	1000	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<0.21	<b>0.63</b>	<0.20	<0.46	NT	NT	NT	<b>0.30</b>	<0.22	<b>1.1</b>	<0.22	
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Naphthalene	mg/kg	4	20	500	<0.21	<0.25	<0.20	<0.46	NT	NT	NT	<0.20	<0.22	<0.24	<0.22	
Phenanthrene	mg/kg	10	500	500	<0.21	<b>2.1</b>	<0.20	<0.46	NT	NT	NT	<b>0.39</b>	<0.22	<b>1.5</b>	<0.22	
Pyrene	mg/kg	1000	1000	1000	<b>0.24</b>	<b>2.3</b>	<0.20	<b>0.56</b>	NT	NT	NT	<b>0.85</b>	<0.22	<b>2.9</b>	<0.22	
<b>VPH</b>																
C5-C8 Aliphatics	mg/kg	100	100	100	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3	
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3	
C9-C10 Aromatics	mg/kg	100	100	100	NT	<6.3	NT	<16	NT	NT	<5.8	NT	<4.9	NT	<4.3	
<b>Target VOCs</b>																
Benzene	mg/kg	2	40	40	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
Ethylbenzene	mg/kg	40	500	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
Naphthalene	mg/kg	4	20	500	NT	<0.31	NT	<0.78	NT	NT	<0.29	NT	<0.24	NT	<0.22	
Toluene	mg/kg	30	500	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
m+p Xylene	mg/kg	100	100	500	NT	<0.13	NT	<0.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
o-Xylene	mg/kg	100	100	500	NT	<0.13	NT	<.31	NT	NT	<0.12	NT	<0.097	NT	<0.087	
<b>Metals</b>																
Antimony	mg/kg	20	20	20	<0.52	<b>2.7</b>	<0.54	<1.1	<0.51	<b>3.7</b>	<0.60	<0.54	<0.57	<b>4.7</b>	<0.55	
Arsenic	mg/kg	20	20	20	<b>3.3</b>	<b>28</b>	<b>3.3</b>	<b>10</b>	<b>3.5</b>	<b>13</b>	<b>3.3</b>	<2.7	<b>3.1</b>	<b>7.3</b>	<b>3.0</b>	
Barium	mg/kg	1000	1000	1000	<b>18</b>	<b>140</b>	<b>27</b>	<b>56</b>	<b>25</b>	<b>250</b>	<b>36</b>	<b>17</b>	<b>27</b>	<b>310</b>	<b>23</b>	
Beryllium	mg/kg	90	90	90	<0.52	<0.68	<0.54	<1.1	<0.51	<0.62	<0.60	<0.54	<0.57	<0.60	<0.55	
Cadmium	mg/kg	70	70	70	<0.52	<b>1.3</b>	<0.54	<1.1	<0.51	<b>4.9</b>	<0.60	<0.54	<0.57	<b>2.0</b>	<0.55	
Chromium (III)	mg/kg	1000	1000	1000	NT	<b>51</b>	NT	<b>41</b>	NT	<b>9.4</b>	NT	NT	<b>16</b>	<b>18</b>	NT	
Chromium (VI)	mg/kg	100	100	100	<b>7.3</b>	<b>3.8</b>	<b>17</b>	<1.00	<b>6.9</b>	<0.49	<b>18</b>	<5.4	<0.47	<0.48	<b>11</b>	
Lead	mg/kg	200	200	200	<b>8.7</b>	<b>230</b>	<b>16</b>	<b>190</b>	<b>22</b>	<b>710</b>	<b>8.4</b>	<b>16</b>	<b>5.5</b>	<b>650</b>	<b>5.0</b>	
Mercury	mg/kg	20	20	20	<0.15	<b>0.58</b>	<0.13	<0.31	<0.15	<b>0.58</b>	<0.17	<0.15	<0.16	<b>2.9</b>	<0.15	
Nickel	mg/kg	600	600	600	<b>7.4</b>	<b>14</b>	<b>23</b>	<b>64</b>	<b>6.3</b>	<b>12</b>	<b>13</b>	<b>5.4</b>	<b>17</b>	<b>18</b>	<b>14</b>	
Selenium	mg/kg	400	400	400	<5.2	<6.8	<5.4	<11	<5.1	<6.2	<6.0	<5.4	<5.7	<6.0	<5.5	
Silver	mg/kg	100	100	100	<2.6	<3.4	<2.7	<5.6	<2.5	<3.1	<3.0	<2.7	<2.8	<3.0	<2.7	
Thallium	mg/kg	8	8	8	<0.52	<0.68	<0.54	<1.1	<0.51	<0.62	<0.60	<0.54	<0.57	<0.60	<0.55	
Vanadium	mg/kg	400	400	400	<b>11</b>	<b>15</b>	<b>20</b>	<b>17</b>	<b>10</b>	<b>10.0</b>	<b>25</b>	<b>8.3</b>	<b>22</b>	<b>19</b>	<b>14</b>	
Zinc	mg/kg	1000	1000	1000	<b>17</b>	<b>220</b>	<b>31</b>	<b>530</b>	<b>19</b>	<b>920</b>	<b>18</b>	<b>16</b>	<b>18</b>	<b>770</b>	<b>69</b>	
<b>VOCs</b>																
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Cis-1,2-Dichloroethylene	mg/kg	0.1	100	100	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	

QC by JMR 5/8/2020

**Abbreviations:**

EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
PAH = Polycyclic Aromatic Hydrocarbons  
NT = Not Tested  
BDL = Below Detection Limit  
MCP = Massachusetts Contingency Plan  
mg/kg = milligram per kilogram

**Notes:**

-- = No standard available  
< = indicates parameter not detected above laboratory method reporting limit, shown  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter equals or exceeds the MCP Method 1, RCS-1 standard (NS Tapley & NWS Bay only)  
**BOLD** Parameter exceeds the most stringent MCP Method 1, S-1 standard  
**BOLD** Parameter exceeds the applicable MCP Method 1, S-2/3 standard  
1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.

**Table 1B**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**Cedar Woods**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)		Sample Location, Depth, and Date														
					RCS-1	S-1/GW-2	S-1/GW-3	SB-106			SB-107		SB-108		SB-109		SB-110		
								0-3 feet	3-4 feet	5-8 feet	0-3 Feet	10-11 Feet	0-3 Feet	6-9 Feet	0-3 Feet	5-8 Feet	DUP-1	0-3 Feet	5-8 Feet
								4/6/2020	4/6/2020	4/6/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020	4/7/2020
<b>EPH</b>																			
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<21	NT	<240	<23	<b>34</b>	<26	<b>92</b>	<28	<78	<83	<22	<110			
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>91</b>	NT	<b>570</b>	<b>36</b>	<b>410</b>	<b>39</b>	<b>1900</b>	<b>42</b>	<b>290</b>	<b>390</b>	<b>68</b>	<b>140</b>			
C11-C22 Aromatics	mg/kg	1000	1000	1000	<b>44</b>	NT	<b>290</b>	<23	<b>76</b>	<0.26	<b>350</b>	<28	<b>100</b>	<b>300</b>	<b>53</b>	<110			
<b>Target PAHS</b>																			
Acenaphthene	mg/kg	4	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Acenaphthylene	mg/kg	1	600	10	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Anthracene	mg/kg	1000	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<b>0.92</b>	<0.22	<1.1			
Benzo(A)Anthracene	mg/kg	7	7	7	<b>0.42</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.79</b>	<0.28	<b>0.78</b>	<b>4</b>	<b>0.59</b>	<1.1			
Benzo(A)Pyrene	mg/kg	2	2	2	<b>0.40</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.70</b>	<0.28	<b>0.91</b>	<b>4.4</b>	<b>0.63</b>	<1.1			
Benzo(B)Fluoranthene	mg/kg	7	7	7	<b>0.35</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.64</b>	<0.28	<b>0.89</b>	<b>4.4</b>	<b>0.57</b>	<1.1			
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<b>0.46</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.51</b>	<0.28	<0.78	<b>3.2</b>	<b>0.42</b>	<1.1			
Benzo(K)Fluoranthene	mg/kg	70	70	70	<b>0.39</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.57</b>	<0.28	<0.78	<b>3.7</b>	<b>0.53</b>	<1.1			
Chrysene	mg/kg	70	70	70	<b>0.52</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.87</b>	<0.28	<b>0.96</b>	<b>5.3</b>	<b>0.74</b>	<1.1			
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<0.24	<0.28	<0.78	<0.83	<0.22	<1.1			
Fluoranthene	mg/kg	1000	1000	1000	<b>0.81</b>	NT	<0.47	<b>0.25</b>	<0.23	<0.26	<b>1.9</b>	<0.28	<b>1.5</b>	<b>9.2</b>	<b>1.3</b>	<1.1			
Fluorene	mg/kg	1000	1000	1000	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.37</b>	<0.28	<0.78	<0.83	<0.22	<1.1			
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<b>0.36</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>0.36</b>	<0.28	<0.78	<b>2.9</b>	<b>0.36</b>	<1.1			
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.57</b>	<0.28	<0.78	<0.83	<0.22	<1.1			
Naphthalene	mg/kg	4	20	500	<0.21	NT	<0.47	<0.23	<0.23	<0.26	<b>0.50</b>	<0.28	<0.78	<b>0.86</b>	<0.22	<1.1			
Phenanthrene	mg/kg	10	500	500	<b>0.44</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>2.1</b>	<0.28	<b>0.80</b>	<b>5.7</b>	<b>1.1</b>	<1.1			
Pyrene	mg/kg	1000	1000	1000	<b>0.88</b>	NT	<0.47	<0.23	<0.23	<0.26	<b>2.1</b>	<0.28	<b>1.7</b>	<b>9</b>	<b>1.4</b>	<1.1			
<b>VPH</b>																			
C5-C8 Aliphatics	mg/kg	100	100	100	NT	NT	<5.5	<b>7.6</b>	<b>26</b>	NT	<b>30</b>	NT	<29	<29	NT	<42			
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	NT	<5.5	<5.4	<b>22</b>	NT	<b>24</b>	NT	<29	<29	NT	<42			
C9-C10 Aromatics	mg/kg	100	100	100	NT	NT	<5.5	<5.4	<b>39</b>	NT	<b>67</b>	NT	<29	<29	NT	<42			
<b>Target VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.30</b>	NT	<0.59	<0.59	NT	<0.83			
Ethylbenzene	mg/kg	40	500	500	NT	NT	<0.11	<0.11	<b>0.48</b>	NT	<b>0.49</b>	NT	<0.59	<0.59	NT	<0.83			
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	NT	<0.11	<0.11	<0.094	NT	<0.11	NT	<0.59	<0.59	NT	<0.83			
Naphthalene	mg/kg	4	20	500	NT	NT	<0.28	<0.27	<b>1.3</b>	NT	<b>1.9</b>	NT	<1.5	<1.5	NT	<2.1			
Toluene	mg/kg	30	500	500	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.42</b>	NT	<0.59	<0.59	NT	<0.83			
m-p Xylene	mg/kg	100	100	500	NT	NT	<0.11	<0.11	<b>1.6</b>	NT	<b>2.7</b>	NT	<0.59	<0.59	NT	<0.83			
o-Xylene	mg/kg	100	100	500	NT	NT	<0.11	<0.11	<0.094	NT	<b>0.45</b>	NT	<0.59	<0.59	NT	<0.83			
<b>Metals</b>																			
Antimony	mg/kg	20	20	20	<b>0.9</b>	NT	<b>0.78</b>	<b>55</b>	<b>5.9</b>	<b>51</b>	<b>10</b>	<b>64</b>	<b>3.5</b>	<b>2.1</b>	NT	<b>3.5</b>			
Arsenic	mg/kg	20	20	20	<b>4</b>	NT	<b>3.2</b>	<b>8.0</b>	<b>3.5</b>	<b>8.1</b>	<b>7.1</b>	<b>23</b>	<b>15</b>	<b>20</b>	<b>2.9</b>	<b>14</b>			
Barium	mg/kg	1000	1000	1000	<b>200</b>	NT	<b>32</b>	<b>280</b>	<b>48</b>	<b>430</b>	<b>470</b>	<b>300</b>	<b>78</b>	<b>86</b>	<b>41</b>	<b>160</b>			
Beryllium	mg/kg	90	90	90	<0.56	NT	<0.59	<0.63	<0.49	<0.65	<0.60	<0.67	<2.0	<2.0	<0.51	<2.8			
Cadmium	mg/kg	70	70	70	<0.56	NT	<0.59	<b>4.4</b>	<b>2.0</b>	<b>5.8</b>	<b>11</b>	<b>3.6</b>	<2.0	<2.0	<0.51	<2.8			
Chromium (III)	mg/kg	1000	1000	1000	NT	NT	NT	<b>33</b>	NT	NT	<b>25</b>	NT	<20	NT	NT	<28			
Chromium (VI)	mg/kg	100	100	100	<b>16</b>	<b>0.50</b>	<b>23</b>	<0.51	<b>19</b>	<b>70</b>	<0.48	<b>64</b>	<1.6	<20	<b>10</b>	<2.2			
Lead	mg/kg	200	200	200	<b>120</b>	NT	<b>150</b>	<b>1200</b>	<b>300</b>	<b>1200</b>	<b>970</b>	<b>1300</b>	<b>180</b>	<b>250</b>	<b>8.1</b>	<b>280</b>			
Mercury	mg/kg	20	20	20	<b>0.17</b>	NT	<0.15	<b>0.60</b>	<0.14	<b>0.18</b>	<b>0.47</b>	<0.19	<0.58	<b>0.99</b>	<0.16	<0.78			
Nickel	mg/kg	600	600	600	<b>12</b>	NT	<b>9.8</b>	<b>76</b>	<b>16</b>	<b>88</b>	<b>40</b>	<b>100</b>	<20	<20	<b>12</b>	<28			
Selenium	mg/kg	400	400	400	<5.36	NT	<5.9	<6.3	<4.9	<6.5	<6.0	<6.7	<20	<20	<5.1	<28			
Silver	mg/kg	100	100	100	<2.8	NT	<3.0	<b>32</b>	<2.5	<b>7.3</b>	<3.0	<b>12</b>	<9.9	<10	<2.5	<14			
Thallium	mg/kg	8	8	8	<0.56	NT	<0.59	<0.63	<0.49	<0.65	<0.60	<0.67	<2.0	<2.0	<0.51	<2.8			
Vanadium	mg/kg	400	400	400	<b>17</b>	NT	<b>12</b>	<b>20</b>	<b>12</b>	<b>10</b>	<b>18</b>	<b>8.1</b>	<20	<20	<b>16</b>	<28			
Zinc	mg/kg	1000	1000	1000	<b>190</b>	NT	<b>91</b>	<b>2900</b>	<b>250</b>	<b>2900</b>	<b>2400</b>	<b>4000</b>	<b>270</b>	<b>250</b>	<b>19</b>	<b>230</b>			
<b>VOCs</b>																			
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	<b>0.19</b>	NT	NT	NT	NT	NT			
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	<b>0.26</b>	NT	NT	NT	NT	NT			
Cis-1,2-Dichloroethylene	mg/kg	0.1	0.1	100	NT	NT	NT	NT	NT	NT	<b>0.20</b>	NT	NT	NT	NT	NT			
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	<b>0.59</b>	NT	NT	NT	NT	NT			
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	<b>0.34</b>	NT	NT	NT	NT	NT			
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	<b>0.30</b>	NT	NT	NT	NT	NT			
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	<b>1.8</b>	NT	NT	NT	NT	NT			
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	<b>0.93</b>	NT	NT	NT	NT	NT			
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	<b>0.46</b>	NT	NT	NT	NT	NT			
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	<b>0.36</b>	NT	NT	NT	NT	NT			
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	<b>6.9</b>	NT	NT	NT	NT	NT			
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	<b>2.3</b>	NT	NT	NT	NT	NT			
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	<b>3.43</b>	NT	NT	NT	NT	NT			

QC by JMR 5/8/2020

**Abbreviations:**  
 EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
 VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
 PAH = Polycyclic Aromatic Hydrocarbons  
 NT = Not Tested  
 BDL = Below Detection Limit  
 MCP = Massachusetts Contingency Plan  
 mg/kg = milligram per kilogram

**Notes:**  
 ~ = No standard available  
 < = indicates parameter not detected  
**BOLD** Parameter detected  
**BOLD** Parameter equals  
**BOLD** Parameter exceeds  
**BOLD** Parameter exceeds  
 1 = Standards are from Massachusetts



**Table 1C**  
**Summary of Soil Analytical Results - April 2020 Soil Borings**  
**Cedar Woods**  
**Framingham, Massachusetts**

Parameter	Units	Reportable Concentrations	Method 1 Cleanup Standards (1)														
			RCS-1	S-1/GW-2	S-1/GW-3	SB-111			SB-112		SB-113			SB-114		SB-115	
						0-1 feet	0-3 feet	4-7 feet	0-3 Feet	5-8 Feet	0-3 feet	5-8 feet	DUP-2	0-3 feet	5-8 feet	0-3 feet	5-8 feet
						4/8/2020	4/8/2020	4/8/2020	4/7/2020	4/7/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020	4/8/2020
<b>EPH</b>																	
C9-C18 Aliphatics	mg/kg	1000	1000	1000	<20	<19	<97	<21	<38	<97	<19	<20	<19	<19	<19	<19	<20
C19-C36 Aliphatics	mg/kg	3000	3000	3000	<b>85</b>	<b>40</b>	<b>340</b>	<b>50</b>	<b>190</b>	<b>450</b>	<19	<20	<b>40</b>	<b>55</b>	<19	<19	<b>57</b>
C11-C22 Aromatics	mg/kg	1000	1000	1000	<b>62</b>	<19	<b>200</b>	<b>22</b>	<b>140</b>	<b>97</b>	<19	<20	<b>220</b>	<b>33</b>	<19	<19	<b>75</b>
<b>Target PAHs</b>																	
Acenaphthene	mg/kg	4	1000	1000	<0.20	<0.19	<b>0.24</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<0.20
Acenaphthylene	mg/kg	1	600	10	<0.20	<0.19	<0.19	<0.21	<0.38	<0.19	<0.19	<0.20	<b>0.77</b>	<0.19	<0.19	<0.19	<0.20
Anthracene	mg/kg	1000	1000	1000	<0.20	<0.19	<b>1.3</b>	<0.21	<b>0.84</b>	<0.19	<0.19	<0.20	<b>0.61</b>	<0.19	<0.19	<0.19	<b>0.33</b>
Benzo(A)Anthracene	mg/kg	7	7	7	<b>1.5</b>	<0.19	<b>4.1</b>	<b>0.39</b>	<b>3</b>	<0.19	<0.19	<0.20	<b>4.9</b>	<b>0.33</b>	<0.19	<0.19	<b>1.6</b>
Benzo(A)Pyrene	mg/kg	2	2	2	<b>2</b>	<0.19	<b>4.1</b>	<b>0.4</b>	<b>3</b>	<0.19	<0.19	<0.20	<b>4.4</b>	<b>0.37</b>	<0.19	<0.19	<b>1.7</b>
Benzo(B)Fluoranthene	mg/kg	7	7	7	<b>2.3</b>	<0.19	<b>3.9</b>	<b>0.37</b>	<b>2.8</b>	<0.19	<0.19	<0.20	<b>3.6</b>	<b>0.29</b>	<0.19	<0.19	<b>1.8</b>
Benzo(G,H,I)Perylene	mg/kg	1000	1000	1000	<b>2</b>	<0.19	<b>3</b>	<b>0.25</b>	<b>2</b>	<0.19	<0.19	<0.20	<b>3.4</b>	<b>0.34</b>	<0.19	<0.19	<b>1.5</b>
Benzo(K)Fluoranthene	mg/kg	70	70	70	<b>1.7</b>	<0.19	<b>3.4</b>	<b>0.38</b>	<b>2.7</b>	<0.19	<0.19	<0.20	<b>3.4</b>	<b>0.31</b>	<0.19	<0.19	<b>1.4</b>
Chrysene	mg/kg	70	70	70	<b>1.8</b>	<0.19	<b>4.4</b>	<b>0.42</b>	<b>3.6</b>	<0.19	<0.19	<0.20	<b>5.7</b>	<b>0.38</b>	<0.19	<0.19	<b>1.9</b>
Dibenz(A,H)Anthracene	mg/kg	0.7	1	1	<b>0.51</b>	<0.19	<b>0.95</b>	<0.21	<b>0.62</b>	<0.19	<0.19	<0.20	<b>0.78</b>	<0.19	<0.19	<0.19	<b>0.38</b>
Fluoranthene	mg/kg	1000	1000	1000	<b>2.2</b>	<0.19	<b>8.5</b>	<b>0.8</b>	<b>6.8</b>	<0.19	<0.19	<0.20	<b>9.3</b>	<b>0.77</b>	<0.19	<0.19	<b>3.3</b>
Fluorene	mg/kg	1000	1000	1000	<0.20	<0.19	<b>0.43</b>	<0.21	<b>0.43</b>	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<b>0.2</b>
Indeno(1,2,3-Cd)Pyrene	mg/kg	7	7	7	<b>1.7</b>	<0.19	<b>2.7</b>	<b>0.21</b>	<b>1.7</b>	<0.19	<0.19	<0.20	<b>2.7</b>	<b>0.25</b>	<0.19	<0.19	<b>1.2</b>
2-Methylnaphthalene	mg/kg	0.7	80	300	<0.20	<0.19	<b>0.22</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<0.20
Naphthalene	mg/kg	4	20	500	<0.20	<0.19	<b>0.35</b>	<0.21	<0.38	<0.19	<0.19	<0.20	<0.19	<0.19	<0.19	<0.19	<b>0.26</b>
Phenanthrene	mg/kg	10	500	500	<b>0.57</b>	<0.19	<b>4.7</b>	<b>0.41</b>	<b>4.7</b>	<0.19	<0.19	<0.20	<b>4.1</b>	<b>0.54</b>	<0.19	<0.19	<b>2.2</b>
Pyrene	mg/kg	1000	1000	1000	<b>1.9</b>	<0.19	<b>6.7</b>	<b>0.73</b>	<b>5.8</b>	<0.19	<0.19	<0.20	<b>11</b>	<b>0.69</b>	<0.19	<0.19	<b>3.5</b>
<b>VPH</b>																	
C5-C8 Aliphatics	mg/kg	100	100	100	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3
C9-C12 Aliphatics	mg/kg	1000	1000	1000	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3
C9-C10 Aromatics	mg/kg	100	100	100	NT	NT	<5.4	NT	<10	NT	<4.5	<4.5	<3.9	NT	NT	NT	<7.3
<b>Target VOCs</b>																	
Benzene	mg/kg	2	40	40	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
Ethylbenzene	mg/kg	40	500	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
Methyl tert-Butyl Ether (MTBE)	mg/kg	0.1	100	100	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
Naphthalene	mg/kg	4	20	500	NT	NT	<b>0.3</b>	NT	<0.52	NT	<0.23	<0.23	<0.20	NT	NT	NT	<0.37
Toluene	mg/kg	30	500	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
m+p Xylene	mg/kg	100	100	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
o-Xylene	mg/kg	100	100	500	NT	NT	<0.11	NT	<0.21	NT	<0.090	<0.090	<0.078	NT	NT	NT	<0.15
<b>Metals</b>																	
Antimony	mg/kg	20	20	20	<b>0.45</b>	<b>29</b>	<b>4.6</b>	<0.50	<b>4.2</b>	<0.56	<0.42	<0.44	<0.48	<b>0.66</b>	<0.40	<0.40	<b>1.7</b>
Arsenic	mg/kg	20	20	20	<1.7	<b>12</b>	<b>8.2</b>	<b>5.0</b>	<b>16</b>	<b>3.7</b>	<b>2.1</b>	<2.2	<2.4	<15	<b>2.6</b>	<b>8.3</b>	
Barium	mg/kg	1000	1000	1000	<b>51</b>	<b>260</b>	<b>80</b>	<b>38</b>	<b>130</b>	<b>48</b>	<b>26</b>	<b>26</b>	<b>17</b>	<b>210</b>	<b>30</b>	<b>94</b>	
Beryllium	mg/kg	90	90	90	<0.34	<0.48	<b>0.44</b>	<0.50	<0.95	<b>0.69</b>	<0.42	<0.044	<0.48	<b>0.59</b>	<b>0.56</b>	<0.64	
Cadmium	mg/kg	70	70	70	<0.34	<b>3.8</b>	<b>0.7</b>	<0.50	<b>1.1</b>	<0.56	<0.42	<0.44	<0.48	<0.59	<0.40	<0.64	
Chromium (III)	mg/kg	1000	1000	1000	NT	NT	<b>13</b>	NT	<b>21</b>	NT	<b>5.7</b>	NT	NT	<29	NT	<b>15</b>	
Chromium (VI)	mg/kg	100	100	100	<b>12</b>	<b>43</b>	<b>2</b>	<b>18</b>	<b>1.3</b>	<b>15</b>	<0.40	<b>6.3</b>	<4.8	<0.59	<b>8</b>	<0.59	
Lead	mg/kg	200	200	200	<b>21</b>	<b>630</b>	<b>460</b>	<b>66</b>	<b>470</b>	<b>36</b>	<b>3.2</b>	<b>3.4</b>	<b>21</b>	<b>270</b>	<b>14</b>	<b>330</b>	
Mercury	mg/kg	20	20	20	<0.14	<b>0.26</b>	<0.16	<0.15	<b>0.80</b>	<0.16	<0.15	<0.14	<0.15	<0.21	<0.13	<0.21	
Nickel	mg/kg	600	600	600	<b>11</b>	<b>52</b>	<b>18</b>	<b>18</b>	<b>17</b>	<5.6	<b>6.2</b>	<b>5.9</b>	<4.8	<29	<b>5.4</b>	<b>9.9</b>	
Selenium	mg/kg	400	400	400	<3.4	<4.8	<4.0	<5.0	<9.5	<5.6	<4.2	<4.4	<4.8	<29	<4.0	<6.4	
Silver	mg/kg	100	100	100	<1.7	<b>4.5</b>	<2.0	<2.5	<4.8	<2.8	<2.1	<2.2	<2.4	<2.9	<2.0	<3.2	
Thallium	mg/kg	8	8	8	<0.34	<b>0.48</b>	<b>0.4</b>	<0.50	<0.95	<0.56	<0.42	<0.44	<0.48	<0.59	<0.40	<0.64	
Vanadium	mg/kg	400	400	400	<b>19</b>	<b>12</b>	<b>14</b>	<b>21</b>	<b>26</b>	<b>16</b>	<b>7.8</b>	<b>8.3</b>	<4.8	<29	<b>11</b>	<b>16</b>	
Zinc	mg/kg	1000	1000	1000	<b>42</b>	<b>1500</b>	<b>330</b>	<b>51</b>	<b>300</b>	<b>43</b>	<b>18</b>	<b>18</b>	<b>40</b>	<b>340</b>	<b>28</b>	<b>210</b>	
<b>VOCs</b>																	
Benzene	mg/kg	2	40	40	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Sec-Butylbenzene	mg/kg	~	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Cis-1,2-Dichloroethylene	mg/kg	0.1	100	100	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Ethylbenzene	mg/kg	40	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Isopropylbenzene (Cumene)	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
P-Isopropyltoluene (p-Cymene)	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Naphthalene	mg/kg	4	20	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
N-Propylbenzene	mg/kg	100	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Toluene	mg/kg	30	500	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Trichloroethylene	mg/kg	0.3	0.3	30	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
1,2,4-Trimethylbenzene	mg/kg	1000	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
1,3,5-Trimethylbenzene	mg/kg	10	~	~	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
Xylenes	mg/kg	100	100	500	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT

QC by JMR 5/8/2020

**Abbreviations:**  
 EPH and VPH= Extractable and Volatile Petroleum Hydrocarbons  
 VOCs and SVOCs = Volatile and Semivolatile Organic Compounds  
 PAH = Polycyclic Aromatic Hydrocarbons  
 NT = Not Tested  
 BDL = Below Detection Limit  
 MCP = Massachusetts Contingency Plan  
 mg/kg = milligram per kilogram

**Notes:**  
 ~ = No standard available  
 < = indicates parameter not detected  
**BOLD** Parameter detected  
**BOLD** Parameter equals  
**BOLD** Parameter exceeds  
**BOLD** Parameter exceeds  
 1 = Standards are from Massachusetts

Table 2  
 Summary of Soil Analytical Results - Shallow Lead Concentrations  
 Cedar Woods  
 618R Waverly Street  
 Framingham, Massachusetts

Parameters	Units	MCP - Method 1 Cleanup Standards <sup>1</sup>	Sample Location, Date, and Depth							
		S-1/GW-3	SB-102	SB-106	SB-107	SB-108	SB-109	SB-110	SB-111	SB-112
			4/6/2020 0-1 feet	4/6/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/7/2020 0-1 feet	4/8/2020 0-1 feet	4/7/2020 0-1 feet
Metals LEAD	mg/Kg	200	73	68	67	900	790	11	21	25

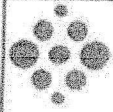
**Abbreviations:**

MCP = Massachusetts Contingency Plan  
 mg/kg = milligram per killogram

**Notes:**

< = indicates parameter not detected above laboratory method reporting limit, shown  
 1 = Standards are from Massachusetts Contingency Plan (MCP), 310 CMR 40, April 2014.  
**BOLD** Parameter detected above laboratory detection limit  
**BOLD** Parameter exceeds the MCP Method 1, S-1/GW-3 Cleanup Standard

**Generator's Waste Profile 1663004-00**



**Stericycle**  
Environmental Solutions

Starts : 07 JUL 2020  
Expires : 06 JUL 2021

Status : PENDING

Sales Rep 1511 Greg Viens  
Acct Mngr 1510 Evan Altmann

**A: GENERATOR ( 646334 ) SITE INFORMATION**

Cedar Woods  
618R Waverly Street  
FRAMINGHAM, MA 01702

EPA EXE  
NAICS 531190 Neshap N  
Phone (508) 532-5470

**B: CUSTOMER ( 39396 ) INFORMATION**

STRATEGIC ENVIRONMENTAL SERVICES, Inc.  
PO BOX 676  
SHREWSBURY, MA 01545-8676

> Contact  
TSDF Approval List No Sub Part P No

**C: WASTE INFORMATION**

On File > MSDS No Analysis Yes Sample No Formulary No

Waste Name SOIL/WATER

Process BORINGS. SOURCE OF SOLVENTS IS UNKNOWN. FORMER AUTO GLASS COMPANY, AUTO PARTS RETAILER, FILLING STATION. SEE ANALYTICAL

Unused Commercial Product No Spill Residue No

**D: PHYSICAL CHARACTERISTICS OF WASTE**

Phys States L-Liq Top Color clear/amber  
Mid Color  
S-Sol Bot Color brown  
% Ash 0  
% Water 50

Odor None  
Layers Bi-Layered  
Spec Grav 1.1  
BTU/Lbs N/A  
% Halogens N/A

PH Range 4-10  
Free Liq % 50  
Flash Test Gen Knowledge  
Flash Rnge >200F  
Viscosity Med  
Pumpable No

**E: CHEMICAL COMPOSITION OF WASTE**

soil/water ( 100 % )  
PCB's 0 Cyanides 0 Phenolics No Sulfides 0 Dioxins No  
TOC <1% VOC <500PPM TAB Profile Information Provided By Generator

**F: METALS METHOD TCLP**

Cadmium <1	Chromium <5	Silver <5	Zinc 0
Arsenic <5	Merc TCLP <0.2	Nickel 0	Copper 0
Barium <100	Lead <5	Thallium 0	Chrome-6

**G: OTHER CHARACTERISTICS OF WASTE**

Ign. Solid No	Oxidizer No	Explosive No	Shock Sensitive No	Cyanide Reactive No	Sulfide Reactive No
Explosive N/A	Asbestos N/A	Radioactive No	Water Reactive No	Reactive (Other) No	
Herbicides No	Pesticides 0	Ammonia No	Infectious No	Medical No	

Gen State UW No

**H: EPA / STATE WASTE IDENTIFICATION**

EPA Waste No	State Waste Yes	TSCA No	Waste Water No	Universal Waste No
Form W113	Source G19	Origin	SubPart CC No	NESHAPS No CERCLA No
			Debris No	Gen State UW No

EPA Codes  
State Codes R015  
UHC

Categorical Discharge Standards No Sub Part P No CTW Category N/A

DW/EHW:

**I: SHIPPING INFORMATION**

Marine Pollutant No

Containers DM Metal Drum

Qty to Ship Now 1

Projected Volume 55/Onetime

DOT Descrip NON-DOT/NON-RCRA REGULATED

**J: SPECIAL DISPOSAL INSTRUCTIONS**

**Generator's Waste Profile 1663004-00**

Status : PENDING



**Stericycle**

Environmental Solutions

Starts : 07 JUL 2020

Expires : 06 JUL 2021

Sales Rep 1511 Greg Viens

Acct Mngr 1510 Evan Altmann

**GENERATOR CERTIFICATION**

To the best of my knowledge and belief, I hereby warrant and represent that the information contained and submitted in this waste profile and all attached documents is true, accurate, and complete and that no material fact has been omitted as to make this misleading. I understand that others may rely on this information in the handling and processing of the waste material described herein. By signing this waste profile, I am certifying that I am authorized to sign such documentation on behalf of the generator.

Thatcher W. Kezer, III

Chief Operating Officer

7/23/2020

Signature

Printed Name

Title

Date

In accordance with 40 CFR 264.12(b), Northland Environmental, LLC has the appropriate permits for, and will accept the waste the generator is shipping as described in this profile.