

# Phase II Environmental Site Assessment

## Location:

5149 County Route 113  
Town of Greenwich, Washinton County, New York 12834

## Prepared for:

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Greenwich, New York 12834

LaBella Project No. 2233938

December 15, 2023



4 British American Boulevard | Latham, NY 12110 | p 518-273-0055

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## 1.0 INTRODUCTION

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LaBella Associates, D.P.C. ("LaBella") was retained by the Town of Greenwich (Client) to conduct a Phase II Environmental Site Assessment (ESA) at the property located at 5149 County Route 113, Town of Greenwich, Washington County, NY hereinafter referred to as the "Site" (see **Figure 1**). This work was conducted following LaBella's August 1, 2023, *Phase I Environmental Site Assessment* for the Site.

The Site is a ( $\pm$ )3.28-acre commercial property that is currently developed with three structures used for storage by Hudson Crossing Park, Inc. LaBella understands that the Site is being redeveloped as a year-round community center with administrative, classroom, and event space, multiple small office buildings, and a construction equipment warehouse/storage yard. Exterior areas of the Site consist of wooded land, maintained yard areas, concrete sidewalks, and a gravel parking lot. The property is bounded by: single-family residential properties to the north, southeast, and south; County Route 113 to the northeast and east; and, the Hudson River to the west.

### 1.1 Limitations & Exceptions

The findings of this Phase II ESA are based on the scope of work and project objectives as stated in LaBella Proposal dated August 18, 2023.

Work associated with this Phase II ESA was performed in accordance with generally accepted environmental engineering and environmental contracting practices for this region. LaBella Associates, D.P.C., makes no other warranty or representation, either expressed or implied, nor is one intended to be included as part of its services, proposals, contracts, or reports.

Phase II ESAs are screening level assessments to assess specific potential environmental concerns identified at a site and are not an exhaustive assessment of environmental conditions on a property. The Phase II ESA is not intended to delineate the nature and extent of contamination at the site, nor address complex geological settings, the fate and transport characteristics of certain hazardous substances, physical limitations imposed by the location of utilities and other man-made objects, and the limitations of assessment technologies.

In addition, LaBella cannot provide guarantees, certifications, or warranties that the property is or is not free of environmental impairment or other regulated solid wastes. The Client shall be aware that the data and representative samples from any given soil sampling point or monitoring well may represent conditions that apply only at that particular location, and such conditions may not necessarily apply to the general Site as a whole.

### 1.2 Background/ Objective

LaBella's August 1, 2023, Phase I ESA for the Site identified the following recognized environmental condition (REC) and significant data gap (SDG). The objective for this Phase II ESA was to investigate Site subsurface soil and groundwater quality relative to the REC and SDG.

1. **REC:** The Subject Property was used as a portion of the former Adirondack School of Northeastern New York between the 1980s and 2010s. According to available information, the structure located at the northwestern portion of the Site was used as classrooms for photograph development, carpentry, stained-glass making, painting & drawing, and print making activities. Chemicals associated with photograph development, unknown waste storage and disposal activities, and discharges to a private septic system from the classroom building may have impacted the environmental integrity of the Site.



2. **SDG:** The Site was developed with the existing buildings in circa 1942. According to Daryl Damas (current occupant), on-site structures have been heated via natural gas, propane, and/or fuel oil and a 275-gallon fuel oil aboveground storage tank (AST) was formerly located near the western corner of the barn building. Additional former heating systems for the on-site buildings are unknown.

This Phase II ESA investigation was completed consistent with the scope of services specified in our proposal (authorized on September 28, 2023) and included the following: a geophysical survey of the subject Property, advancement of three (3) soil borings; and collection of three (3) soil samples submitted for laboratory analysis. Labella's methods, observations, results, conclusions, and recommendations are included in the following sections. Supporting documentation is attached.

## **2.0 FIELD INVESTIGATION**

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### **2.1 Geophysical Survey and Results**

On October 16, 2023, LaBella conducted a geophysical survey (subsurface utility mark out) using ground penetrating radar (GPR) and magnetometry technologies to explore for anomalies indicative of potential underground storage tanks (USTs) or backfilled excavations and to locate nearby subsurface utilities and structures to aid in the safe placement of soil borings.

One anomaly suggestive of a potential UST was identified along the northwest perimeter of the barn building. Detailed results of the geophysical survey are shown on **Figure 2**.

Note: LaBella makes an attempt to locate the underground utilities and subsurface structures in the investigation area; however, not all utilities and structures are detectable. It should be understood that the only way to confirm the identity of a detected utility is to expose the feature through careful excavation. Caution should be taken when drilling or excavating.

### **2.2 Subsurface Soil Investigation**

Prior to the initiation of subsurface work, LaBella submitted a location request to *UDig New York* to locate utilities in the areas near the property boundaries.

#### **2.2.1 Soil Borings**

On October 24, 2023, Core Down Drilling of Brewster, New York, advanced three (3) soil borings at the Site under direct supervision of a LaBella environmental professional. Soil borings SB-01 through SB-03 were extended at downgradient locations relative to the classroom building, the septic system's leach field, and near the subsurface anomaly, respectively. Soil boring locations are depicted on **Figure 3**.

Soil borings were extended using a track-mounted Geoprobe® Systems Model 54DT direct-push sampling system to 16-feet (ft) below ground surface (bgs) and were strategically placed at a safe distance away from the existing subsurface features. The use of direct-push technology allowed for continuous sampling, observation, and characterization of overburden soils. Samples were collected using a MacroCore® sampler equipped with disposable polyethylene sleeves. Soil cores were retrieved in 4-foot sections and are cut from the polyethylene sleeves for observation and sampling. The MacroCore® sampler was decontaminated between boring locations using an Alconox detergent and potable water solution followed by a potable water rinse.



Soils extracted from borings were continuously assessed from the ground surface to the bottom of each boring. LaBella personnel visually and physically examined soil samples and recorded observations of the general lithology, visible layering, evidence of non-native fill/historic fill materials, indications of chemical or other staining, odors, and other distinctive features. Portions of the soil from borings were field screened for the presence of volatile organic compounds (VOCs) using a photoionization detector (PID) equipped with a 10.6 electronvolt (eV) lamp. Groundwater was not encountered during this investigation. Soil boring logs were completed for each soil boring and are included in **Appendix 1**.

### 2.2.2 *Soil Sampling*

Three (3) soil samples (one from each boring/investigation area) were selected for the laboratory analyses. The table below summarizes sample identifications, areas/locations, depths, laboratory analyses. Soil samples were submitted for analysis under standard Chain of Custody procedures to Phoenix Environmental Laboratories (Phoenix) of Manchester, Connecticut, a New York State Department of Health (NYSDOH) Environmental Laboratory Accreditation Program (ELAP) certified laboratory.

**Table A: Sample Location and Analysis Summary**

Sample IDs	Exploration Area and Location	Soil Sample Depth (ft bgs)	Laboratory Analyses
SB-01 (6-7ft)	REC: Downgradient of former classroom building (with historical photograph developing activities and unknown waste storage/disposal methods) and the septic system leach field.	6 to 7	TCL VOCs, TCL SVOCS, TAL Metals
SB-02 (8-10ft)		8 to 10	
SB-03 (14-15ft)	SDG: Downgradient of anomaly adjacent to the barn building. Former heating systems of the on-site buildings are unknown.	14 to 15	CP-51 VOCs, CP-51 SVOCS

- New York State Department of Environmental Conservation (NYSDEC) Part 375 Target Compound List (TCL) VOCs and NYSDEC Commissioner Policy 51 (CP-51) list VOCs via USEPA Method 8260,
- NYSDEC Part 375 TCL semi-volatile organic compounds (SVOCS) and CP-51 list SVOCS via USEPA Method 8270
- Target Analyte List (TAL) metals via USEPA Methods 6010 and 7471.

## 3.0 FINDINGS

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### 3.1 Site Geology and Hydrology

Soil borings were advanced to a maximum depth of 16 feet bgs. Non-native materials encountered at the Site consisted of topsoil to approximately 0.5 feet bgs at the three soil borings. Soils at the Site generally consisted of silty sands with fine gravel, underlain by gravel-sand-silt mixtures. Bedrock and groundwater were not encountered in the borings. Groundwater is assumed to flow in a southern/southwestern direction, towards the Hudson River. Soil Boring Logs are included as **Appendix 1** and the soil boring locations are shown on **Figure 3**.

### 3.2 Field Screening Results

No visual or olfactory evidence of impairment including petroleum odors or stained soil were observed in soil borings SB-01 through SB-03. PID measurements were less than 1 part per million (ppm) in three soil borings.

### 3.3 Laboratory Analytical Results

Laboratory data was reviewed, tabulated, and compared to applicable NYSDEC standards, criteria, and guidance (SCGs). An analytical results summary table is attached (**Table 1**). Detected



concentrations of VOCs and SVOCs, and TAL metals were compared to 6 NYCRR Part 375-6.8 Unrestricted Use Soil Cleanup Objectives (UUSCOs). Although applicable to sites in a NYSDEC remediation program, the soil results were also compared to the Part 375 Residential Use SCOs (RUSCOs) to provide an additional point of reference for property evaluations. The laboratory report is attached as **Appendix 2**.

- No TCL VOCs or SVOCs were detected above laboratory method detection limits (MDLs) in SB-01(6-7ft) or SB-02(7-8ft) and no CP-51 VOCs or SVOCs were detected above the MDLs in SB-03 (14-15ft). No PID readings greater than 1.0 part per million were recorded during fieldwork activities.
- Nickel and zinc were detected at concentrations exceeding their UUSCOs in soil boring SB-02(8-10ft); however, their detected concentrations did not exceed corresponding RUSCOs. Other TAL metals were detected at trace levels (less than the UUSCOs) or were not detected above laboratory method detection limits (MDLs).

## 4.0 CONCLUSIONS AND RECOMMENDATIONS

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LaBella was retained by the Town of Greenwich to conduct a Phase II ESA at the Subject Property located at 5149 County Route 113, Town of Greenwich, Washington County, New York. The Phase II ESA consisted of a geophysical survey of the Subject Property, advancement of three (3) soil borings; and collection of three (3) soil samples submitted for laboratory analyses.

The geophysical survey identified evidence of one (1) anomaly suggestive of a potential UST east of the barn building. The septic system, subsurface utilities (water, electrical, communications, and gas lines) were also identified.

No visual or olfactory evidence of a chemical or petroleum release were observed in the three soil borings, and no VOCs or SVOCs were detected above laboratory method detection limits in corresponding soil samples SB-01(6-7ft), SB-02(7-8ft), and SB-03(14-15ft). Analytical results for metals at boring SB-01 were either non-detect or below UUSCOs. Although zinc and nickel concentrations at SB-02 exceeded their respective UUSCOs, these concentrations did not exceed their respective RUSCOs do not require remediation. Based on the collected data, the former Site use as the Adirondack School of Northeastern New York (REC) is no longer considered a REC. Potential future Site redevelopment that disturbs the Subject property soil should ensure that material is properly managed. We note that these metals mean that the soil does not meet the definition of clean fill and it cannot be transported off-Site for unrestricted use.

This investigation resolved the SDG and did not identify a REC associated with the anomaly near the barn. However, LaBella recommends that this anomaly be uncovered (i.e., daylighted) to determine if it is an underground storage tank (UST) for petroleum. If a UST is confirmed, it should be removed consistent with NYSDEC criteria.

If different subsurface conditions are discovered during Site redevelopment (e.g., impacted soil, debris, tank, etc.), that material should also be handled and disposed of properly, with reporting to the NYSDEC as warranted.



## 5.0 SIGNATURE OF ENVIRONMENTAL PROFESSIONAL

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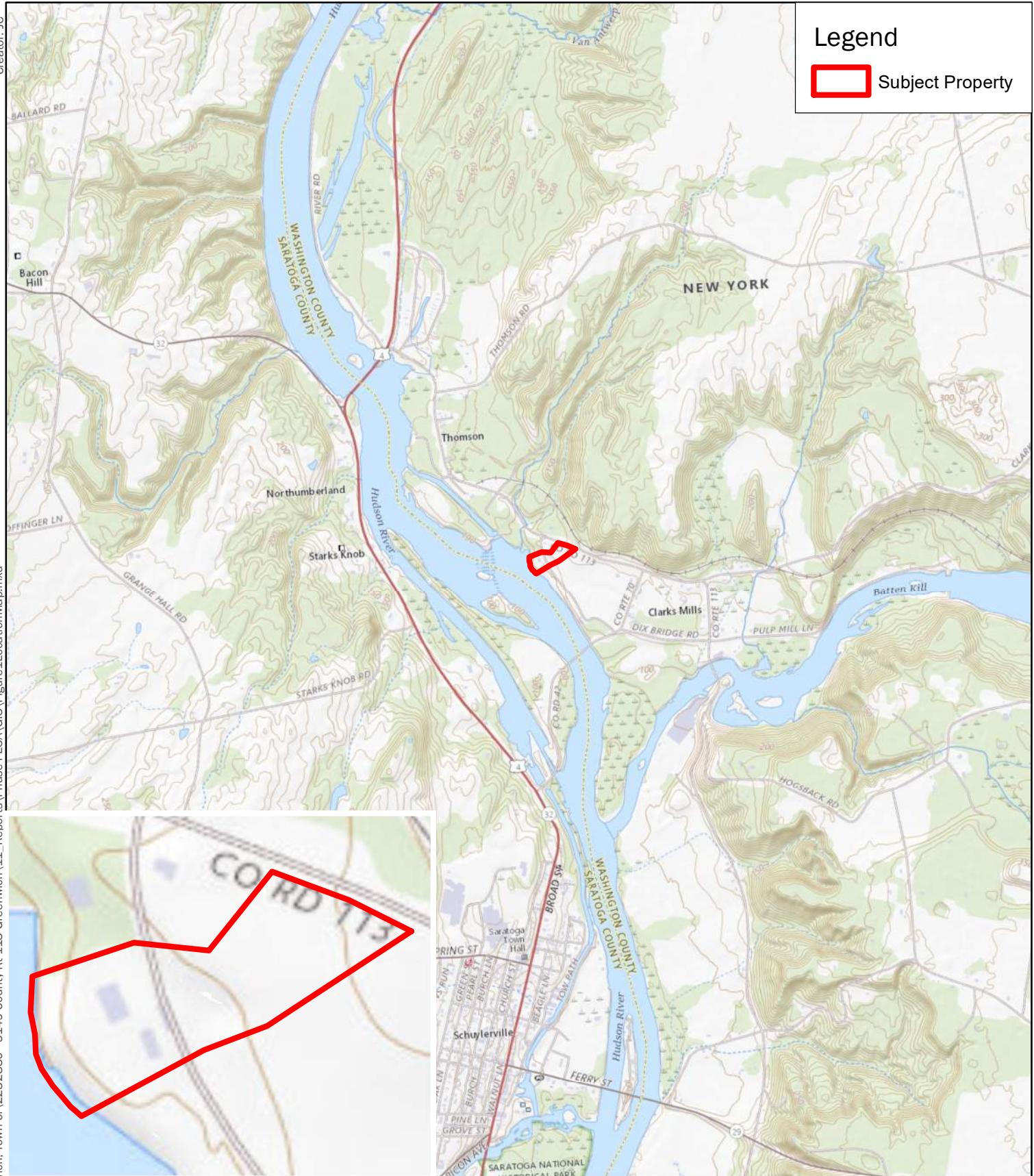
Report Prepared By:

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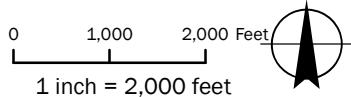
Branson Fields – LaBella Associates  
Environmental Scientist – Project Manager



## FIGURES



Source:  
Washington County 2023 Tax Parcel Dataset;  
US Topo. 2018 - USGSTopo (MapServer)  
Layer: USGS TNM Topo Base Map. Accessed  
19 July 2023



Town  
of Greenwich

Town of Greenwich Hudson  
Riverside Park  
5149 County Route 113  
Town of Greenwich  
Washington County, NY  
LaBella Project No: 2232830  
Date: 8/1/2023

**Location Map**

**FIGURE #1**

# LEGEND

ALL UTILITIES ARE INDICATED WITH DOTTED AND DASHED LINES. PLEASE SEE BELOW COLOR CODE FOR LINE DESIGNATION.

## UTILITY COLOR CODE

<span style="color: red;">█</span>	ELECTRIC
<span style="color: yellow;">█</span>	GAS ( INCLUDES ALL GASSES LIQUID OR NON LIQUID)
<span style="color: blue;">█</span>	POTABLE WATER
<span style="color: green;">█</span>	SEWER/ DRAINAGE
<span style="color: magenta;">█</span>	TEMPORARY SURVEY
<span style="border: 1px solid black; width: 1em; height: 1em;"></span>	(WHITE) DEFINED WORK ZONE
<span style="color: purple;">█</span>	NON POTABLE WATER/ IRRIGATION/ SLURRY LINES
<span style="color: orange;">█</span>	COMMUNICATIONS/ ALARM/ SIGNAL
	LaBella Associates

Figure 2 - Geophysical Survey

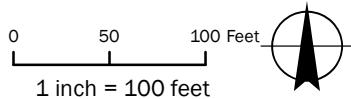


## Legend

- Subject Property
- + Soil Boring Location
- Septic Leach Field
- Subsurface Anomaly Location



Source:  
Washington County 2023 Tax Parcel Dataset;  
NYS Department of Transportation 2022  
Simplified Street Dataset; ARCGIS Basemap  
Layer Orthophotograph dated March 2021.



Town  
of Greenwich

Town of Greenwich Hudson  
Riverside Park  
5149 County Route 113  
Town of Greenwich  
Washington County, NY  
LaBella Project No: 2233938  
Date: 11/30/2023

**Soil Boring  
Location Map**  
**FIGURE #3**



# TABLE

**Table 1**  
 Soil Sample Laboratory Analytical Results Summary  
 5149 County Route 113  
 Town of Greenwich, New York  
 LaBella Project No. 2233938

Sample Location	NYSDEC Part 375-6 Soil Cleanup Objectives			Downgradient "Classroom" Building	Downgradient Septic Leach Field	Downgradient Anomaly, "Barn" Building		
LaBella Sample ID (depth)	CP-51 Soil - Gas/Fuel Oil	Unrestricted Use	Residential Use	SB-01 (6-7FT)	SB-02 (7-8FT)	SB-03 (14-15FT)		
Lab Sample ID				CP33746	CP33747	CP33748		
Collection Date				10/24/2023	10/24/2023	10/24/2023		
Matrix				Soil	Soil	Soil		
Compound	Results							
<b>Miscellaneous/Inorganics</b>	<b>%</b>			<b>%</b>				
Percent Solid	~	~	~	77	87	89		
<b>Metals, Total</b>	<b>mg/kg</b>			<b>mg/kg</b>				
Aluminum	~	~	~	17,100	32,600	NT		
Antimony	~			ND	ND	NT		
Arsenic	~	13	16	4.08	4.93	NT		
Barium	~	350	350	122	90.5	NT		
Beryllium	~	7.2	14	0.82	0.97	NT		
Cadmium	~	2.5	2.5	0.91	1.99	NT		
Calcium	~	~	~	4,900	1,380	NT		
Chromium	~	30	30	18.3	10.8	NT		
Cobalt	~	~	~	11.4	15	NT		
Copper	~	50	270	15.1	36.6	NT		
Iron	~			33,600	69,100	NT		
Lead	~	63	400	9	11.6	NT		
Magnesium	~			4,260	11,200	NT		
Manganese	~	1,600	2,000	340	1,140	NT		
Mercury	~	0.18	0.81	ND	ND	NT		
Nickel	~	30	140	17.2	55.9	NT		
Potassium	~	~	~	1,390	620	NT		
Selenium	~	3.9	36	ND	ND	NT		
Silver	~	2	36	ND	ND	NT		
Sodium	~	~	~	263	28.2	NT		
Thallium	~	~	~	ND	ND	NT		
Vanadium	~	~	~	36.6	23.2	NT		
Zinc	~	109	2,200	73.5	174	NT		
<b>Volatiles- STARS/CP-51 By SW8260D</b>	<b>mg/kg</b>			<b>mg/kg</b>				
1,2,4-Trimethylbenzene	3.6	3.6	47	NT	NT	ND		
1,3,5-Trimethylbenzene	8.4	8.4	47	NT	NT	ND		
Benzene	0.06	0.06	2.9	NT	NT	ND		
Ethylbenzene	1	1	30	NT	NT	ND		
Isopropylbenzene	2.3	~	~	NT	NT	ND		
m&p-Xylene	~	~	~	NT	NT	ND		
Methyl t-Butyl Ether (MTBE)	0.93	0.93	62	NT	NT	ND		
Naphthalene	12	12	100	NT	NT	ND		
n-Butylbenzene	12	12	100	NT	NT	ND		
n-Propylbenzene	3.9	3.9	100	NT	NT	ND		
o-Xylene	~	~	~	NT	NT	ND		
p-Isopropyltoluene	10	~	~	NT	NT	ND		
sec-Butylbenzene	11	11	100	NT	NT	ND		
tert-Butylbenzene	5.9	5.9	100	NT	NT	ND		
Toluene	0.7	0.7	100	NT	NT	ND		
Total Xylenes	0.26	0.26		NT	NT	ND		
<b>Volatiles (TCL) By SW8260D</b>	<b>mg/kg</b>			<b>mg/kg</b>				
1,1,1-Trichloroethane	~	0.68	100	ND	ND	NT		
1,1,2,2-Tetrachloroethane	~	~	~	ND	ND	NT		
1,1,2-Trichloroethane	~	~	~	ND	ND	NT		
1,1-Dichloroethane	~	0.27	19	ND	ND	NT		
1,1-Dichloroethene	~	0.33	100	ND	ND	NT		
1,2,3-Trichlorobenzene	~	~	~	ND	ND	NT		
1,2,4-Trichlorobenzene	~	~	~	ND	ND	NT		
1,2-Dibromo-3-chloropropane	~	~	~	ND	ND	NT		
1,2-Dibromoethane	~	~	~	ND	ND	NT		
1,2-Dichlorobenzene	~	1.1	100	ND	ND	NT		
1,2-Dichloroethane	~	0.02	2.3	ND	ND	NT		
1,2-Dichloropropane	~	~	~	ND	ND	NT		
1,3-Dichlorobenzene	~	2.4	17	ND	ND	NT		
1,4-Dichlorobenzene	~	1.8	9.8	ND	ND	NT		
2-Hexanone	~	~	~	ND	ND	NT		
4-Methyl-2-pentanone	~	~	~	ND	ND	NT		
Acetone	~	0.05	100	ND	ND	NT		
Benzene	0.06	0.06	2.9	ND	ND	NT		
Bromochloromethane	~	~	~	ND	ND	NT		
Bromodichloromethane	~	~	~	ND	ND	NT		
Bromoform	~	~	~	ND	ND	NT		
Bromomethane	~	~	~	ND	ND	NT		
Carbon Disulfide	~	~	~	ND	ND	NT		
Carbon tetrachloride	~	0.76	1.4	ND	ND	NT		
Chlorobenzene	~	1.1	100	ND	ND	NT		
Chloroethane	~	~	~	ND	ND	NT		
Chloroform	~	0.37	10	ND	ND	NT		
Chloromethane	~	~	~	ND	ND	NT		
cis-1,2-Dichloroethene	~	0.25	59	ND	ND	NT		
cis-1,3-Dichloropropene	~	~	~	ND	ND	NT		
Cyclohexane	~	~	~	ND	ND	NT		
Dibromochloromethane	~	~	~	ND	ND	NT		
Dichlorodifluoromethane	~	~	~	ND	ND	NT		
Ethylbenzene	1	1	30	ND	ND	NT		
Isopropylbenzene	2.3	~	~	ND	ND	NT		
m&p-Xylene	~	~	~	ND	ND	NT		
Methyl ethyl ketone	~	0.12	100	ND	ND	NT		
Methyl t-butyl ether (MTBE)	0.93	0.93	62	ND	ND	NT		
Methylacetate	~	~	~	ND	ND	NT		
Methylcyclohexane	~	~	~	ND	ND	NT		
Methylene chloride	~	0.05	51	ND	ND	NT		
o-Xylene	~	~	~	ND	ND	NT		
Styrene	~	~	~	ND	ND	NT		
Tetrachloroethene	~	1.3	5.5	ND	ND	NT		
Toluene	0.7	0.7	100	ND	ND	NT		
Total Xylenes	0.26	0.26	~	ND	ND	NT		
trans-1,2-Dichloroethene	~	0.19	100	ND	ND	NT		
trans-1,3-Dichloropropene	~	~	~	ND	ND	NT		
Trichloroethene	~	0.47	10	ND	ND	NT		
Trichlorofluoromethane	~	~	~	ND	ND	NT		
Trichlorotrifluoroethane	~	~	~	ND	ND	NT		
Vinyl chloride	~	0.02	0.21	ND	ND	NT		

**Table 1**  
 Soil Sample Laboratory Analytical Results Summary  
 5149 County Route 113  
 Town of Greenwich, New York  
 LaBella Project No. 2233938

Sample Location	NYSDEC Part 375-6 Soil Cleanup Objectives			Downgradient "Classroom" Building	Downgradient Septic Leach Field	Downgradient Anomaly, "Barn" Building		
LaBella Sample ID (depth)	CP-51 Soil - Gas/Fuel Oil	Unrestricted Use	Residential Use	SB-01 (6-7FT)	SB-02 (7-8FT)	SB-03 (14-15FT)		
Lab Sample ID				CP33746	CP33747	CP33748		
Collection Date				10/24/2023	10/24/2023	10/24/2023		
Matrix				Soil	Soil	Soil		
Compound				Results				
<b>Semivolatiles-STARS/CP-51 By SW8270D</b>	<b>mg/kg</b>			<b>mg/kg</b>				
Acenaphthene	20	20	100	NT	NT	ND		
Acenaphthylene	100	100	100	NT	NT	ND		
Anthracene	100	100	100	NT	NT	ND		
Benz(a)anthracene	1	1	1	NT	NT	ND		
Benzo(a)pyrene	1	1	1	NT	NT	ND		
Benzo(b)fluoranthene	1	1	1	NT	NT	ND		
Benzo(ghi)perylene	100	100	100	NT	NT	ND		
Benzo(k)fluoranthene	0.8	0.8	1	NT	NT	ND		
Chrysene	1	1	1	NT	NT	ND		
Dibenz(a,h)anthracene	0.33	0.33	0.33	NT	NT	ND		
Fluoranthene	100	100	100	NT	NT	ND		
Fluorene	30	30	100	NT	NT	ND		
Indeno(1,2,3-cd)pyrene	0.5	0.5	0.5	NT	NT	ND		
Naphthalene	12	12	100	NT	NT	ND		
Phenanthrene	100	100	100	NT	NT	ND		
Pyrene	100	100	100	NT	NT	ND		
<b>Semivolatiles By SW8270D</b>	<b>mg/kg</b>			<b>mg/kg</b>				
1,1-Biphenyl	~	~	~	ND	ND	NT		
1,2,4,5-Tetrachlorobenzene	~	~	~	ND	ND	NT		
2,2'-Oxybis(1-Chloropropane)	~	~	~	ND	ND	NT		
2,3,4,6-tetrachlorophenol	~	~	~	ND	ND	NT		
2,4,5-Trichlorophenol	~	~	~	ND	ND	NT		
2,4,6-Trichlorophenol	~	~	~	ND	ND	NT		
2,4-Dichlorophenol	~	~	~	ND	ND	NT		
2,4-Dimethylphenol	~	~	~	ND	ND	NT		
2,4-Dinitrophenol	~	~	~	ND	ND	NT		
2,4-Dinitrotoluene	~	~	~	ND	ND	NT		
2,6-Dinitrotoluene	~	~	~	ND	ND	NT		
2-Chloronaphthalene	~	~	~	ND	ND	NT		
2-Chlorophenol	~	~	~	ND	ND	NT		
2-Methylnaphthalene	~	~	~	ND	ND	NT		
2-Methylphenol (o-cresol)	~	0.33	100	ND	ND	NT		
2-Nitroaniline	~	~	~	ND	ND	NT		
2-Nitrophenol	~	~	~	ND	ND	NT		
3&4-Methylphenol (m&p-cresol)	~	~	~	ND	ND	NT		
3,3'-Dichlorobenzidine	~	~	~	ND	ND	NT		
3-Nitroaniline	~	~	~	ND	ND	NT		
4,6-Dinitro-2-methylphenol	~	~	~	ND	ND	NT		
4-Bromophenyl phenyl ether	~	~	~	ND	ND	NT		
4-Chloro-3-methylphenol	~	~	~	ND	ND	NT		
4-Chloroaniline	~	~	~	ND	ND	NT		
4-Chlorophenyl phenyl ether	~	~	~	ND	ND	NT		
4-Nitroaniline	~	~	~	ND	ND	NT		
4-Nitrophenol	~	~	~	ND	ND	NT		
Acenaphthene	20	20	100	ND	ND	NT		
Acenaphthylene	100	100	100	ND	ND	NT		
Acetophenone	~	~	~	ND	ND	NT		
Anthracene	100	100	100	ND	ND	NT		
Atrazine	~	~	~	ND	ND	NT		
Benz(a)anthracene	1	1	1	ND	ND	NT		
Benzaldehyde	~	~	~	ND	ND	NT		
Benzo(a)pyrene	1	1	1	ND	ND	NT		
Benzo(b)fluoranthene	1	1	1	ND	ND	NT		
Benzo(ghi)perylene	100	100	100	ND	ND	NT		
Benzo(k)fluoranthene	0.8	0.8	1	ND	ND	NT		
Benzyl butyl phthalate	~	~	~	ND	ND	NT		
Bis(2-chloroethoxy)methane	~	~	~	ND	ND	NT		
Bis(2-chloroethyl)ether	~	~	~	ND	ND	NT		
Bis(2-ethylhexyl)phthalate	~	~	~	ND	ND	NT		
Caprolactam	~	~	~	ND	ND	NT		
Carbazole	~	~	~	ND	ND	NT		
Chrysene	1	1	1	ND	ND	NT		
Dibenz(a,h)anthracene	0.33	0.33	0.33	ND	ND	NT		
Dibenzofuran	~	7	14	ND	ND	NT		
Diethyl phthalate	~	~	~	ND	ND	NT		
Dimethylphthalate	~	~	~	ND	ND	NT		
Di-n-butylphthalate	~	~	~	ND	ND	NT		
Di-n-octylphthalate	~	~	~	ND	ND	NT		
Fluoranthene	100	100	100	ND	ND	NT		
Fluorene	30	30	100	ND	ND	NT		
Hexachlorobenzene	~	0.33	0.33	ND	ND	NT		
Hexachlorobutadiene	~	~	~	ND	ND	NT		
Hexachlorocyclopentadiene	~	~	~	ND	ND	NT		
Hexachloroethane	~	~	~	ND	ND	NT		
Indeno(1,2,3-cd)pyrene	0.5	0.5	0.5	ND	ND	NT		
Isophorone	~	~	~	ND	ND	NT		
Naphthalene	12	12	100	ND	ND	NT		
Nitrobenzene	~	~	~	ND	ND	NT		
N-Nitrosodimethylamine	~	~	~	ND	ND	NT		
N-Nitrosodi-n-propylamine	~	~	~	ND	ND	NT		
N-Nitrosodiphenylamine	~	~	~	ND	ND	NT		
Pentachlorophenol	~	0.8	2.4	ND	ND	NT		
Phenanthrene	100	100	100	ND	ND	NT		
Phenol	~	0.33	100	ND	ND	NT		
Pyrene	100	100	100	ND	ND	NT		
<b>1,4-dioxane By SW8260D</b>	<b>mg/kg</b>			<b>mg/kg</b>				
1,4-dioxane	~	0.1	9.8	ND	ND	NT		

**Notes:**

Exceedances of NYSDEC Part 375-6 soil cleanup objectives (SCOs) are formatted consistent with the SCO column headers.

mg/kg= milligrams per kilogram or parts per million (ppm)

% = percentage

ND - analyte not detected at or above the level indicated.

NT - Indicates the analyte was not a target for this sample

~ - Indicates that no regulatory limit has been established for this analyte.



# APPENDIX 1

**Soil Boring Logs**

 <b>LaBella</b> Powered by partnership. Soil Boring Log 4 British American Boulevard, Latham, NY			<b>PROJECT</b> Phase II Environmental Site Assessment 5149 County Route 113 Town of Greenwich, Washington County, New York			<b>BORING:</b> SB - 01 <b>SHEET:</b> 1 of 1 <b>JOB:</b> 2233938 <b>CHKD BY:</b> BF <b>DATE:</b> 10/24/2023		
CONTRACTOR: Core Down Drilling DRILLER: Joe Bellucci LABELLA REPRESENTATIVE: Branson Fields			BORING LOCATION: Rear/Downgradient of "Classroom" Building GROUND SURFACE ELEVATION NA START DATE: 10/24/23 END DATE: 10/24/23			TIME: 09:15 TO 09:45 DATUM: NA WEATHER: M.Cloudy, ~40deg.F		
TYPE OF DRILL RIG: Geoprobe 54DT AUGER SIZE AND TYPE: NA OVERBURDEN SAMPLING METHOD: Direct Push								
DEPTH (FEET BGS)	SAMPLE			VISUAL CLASSIFICATION			PID FIELD SCREEN (PPM)	REMARKS
	SAMPLE RECOVERY (INCHES)	SAMPLE NO. AND DEPTH	STRATA CHANGE (FEET BGS)					
0	36/48	No. 01 0 to 4-ft. bgs	Topsoil	6" Topsoil with vegetation			<1	Soil Sample: SB-01(6to7ft) 9:55
1			SM	30" Very loose, light brown, silty sand (vf-m), dry, non-plastic, NOSOI.				
2				12" Same as above, moist, NOSOI.				
3				12" Very loose, light brown, silty sand (f-c) with gravel (A), moist, non-plastic, NOSOI.				
4	30/48	No. 02 4 to 8-ft. bgs	SM	6" Same as above, NOSOI.			<1	
5				12" Same as above, moist, NOSOI.				
6				12" Very loose, light brown, silty sand (f-c) with gravel (A), moist, non-plastic, NOSOI.				
7				24" Very loose, light brown, silty gravel (SA) with sand (f-c), moist, non-plastic, NOSOI.				
8	24/48	No. 03 8 to 12 ft. bgs.	GM	24" Same as above, NOSOI			<1	
9				24" Same as above, NOSOI				
10				24" Same as above, NOSOI				
11				24" Same as above, NOSOI				
12	24/48	No. 04 12 to 16-ft. bgs	GM	24" Same as above, NOSOI			<1	
13				24" Same as above, NOSOI				
14				24" Same as above, NOSOI				
15				24" Same as above, NOSOI				
16				End of Boring at 16-ft. bgs.				
17								
18								
19								
20								
21								
22								
23								
24								
25								
			DEPTH (FT)			NOTES:		
WATER LEVEL DATA			BOTTOM OF CASING	BOTTOM OF BORING	GROUNDWATER ENCOUNTERED			
DATE	TIME	ELAPSED TIME			na			
GENERAL NOTES								
1) STRATIFICATION LINES REPRESENT APPROXIMATE BOUNDARY BETWEEN SOIL TYPES, TRANSITIONS MAY BE GRADUAL.								
2) WATER LEVEL READINGS HAVE BEEN MADE AT TIMES AND UNDER CONDITIONS STATED, FLUCTUATIONS OF GROUNDWATER								
BGS = Below Ground Surface NA = Not Applicable			and = 35 - 50% some = 20 - 35% little = 10 - 20% trace = 1 - 10%			C = Coarse M = Medium F = Fine VF = Very Fine		
						R = Rounded A = Angular SR = Subrounded SA = Subangular		
BORING: SB - 01								

 <b>LaBella</b> Powered by partnership. Soil Boring Log 4 British American Boulevard, Latham, NY			<b>PROJECT</b> Phase II Environmental Site Assessment 5149 County Route 113 Town of Greenwich, Washington County, New York			<b>BORING:</b> SB - 02 <b>SHEET:</b> 1 of 1 <b>JOB:</b> 2233938 <b>CHKD BY:</b> BF <b>DATE:</b> 10/24/2023	
CONTRACTOR: Core Down Drilling DRILLER: Joe Bellucci LABELLA REPRESENTATIVE: Branson Fields			BORING LOCATION: South/Downgradient of Septic Leach Field GROUND SURFACE ELEVATION NA START DATE: 10/24/23 END DATE: 10/24/23			TIME: 10:00 TO 10:30 DATUM: NA WEATHER: M.Cloudy, ~40deg.F	
TYPE OF DRILL RIG: Geoprobe 54DT AUGER SIZE AND TYPE: NA OVERBURDEN SAMPLING METHOD: Direct Push						DRIVE SAMPLER TYPE: Macrocore INSIDE DIAMETER: 2-inch OTHER: 4-feet core intervals	
DEPTH (FEET BGS)	SAMPLE			VISUAL CLASSIFICATION		PID FIELD SCREEN (PPM)	REMARKS
	SAMPLE RECOVERY (INCHES)	SAMPLE NO. AND DEPTH	STRATA CHANGE (FEET BGS)				
0	39/48	No. 01 0 to 4-ft. bgs	Topsoil	6" Topsoil with vegetation		<1	Soil Sample: SB-02(7-8ft) 10:45
1			SM	33" Very loose, light brown, silty sand (vf-m), dry, non-plastic, NOSOI.			
2				21" Same as above, moist, NOSOI.			
3				9" Very loose, light brown, silty sand (f-c) with gravel (A); moist, non-plastic, NOSOI.			
4	30/48	No. 02 4 to 8-ft. bgs	SM	25" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.		<1	<1
5			SM	24" Very loose, red/orange, silty gravel (SR), moist, non-plastic, little sand (f-m), NOSOI			
6				22" Very loose, light brown, silty sand (f-c) with gravel (A); moist, non-plastic, NOSOI.			
7				10" Very loose, light brown, silty sand (f-c) with gravel (A); moist, non-plastic, NOSOI.			
8	25/48	No. 03 8 to 12 ft. bgs.	GM	18" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.		<1	<1
9			GM	16" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
10				14" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
11				12" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
12	24/48	No. 04 12 to 16-ft. bgs	GM	10" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.		<1	<1
13			GM	8" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
14				6" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
15				4" Very loose, dark brown, silty gravel (SR) with sand (f-m), moist, non-plastic, NOSOI.			
16				End of Boring at 16-ft. bgs.			
WATER LEVEL DATA			BOTTOM OF CASING	BOTTOM OF BORING	GROUNDWATER ENCOUNTERED	NOTES:	
DATE	TIME	ELAPSED TIME					
					na		
GENERAL NOTES							
1) STRATIFICATION LINES REPRESENT APPROXIMATE BOUNDARY BETWEEN SOIL TYPES, TRANSITIONS MAY BE GRADUAL.							
2) WATER LEVEL READINGS HAVE BEEN MADE AT TIMES AND UNDER CONDITIONS STATED, FLUCTUATIONS OF GROUNDWATER							
BGS = Below Ground Surface		and = 35 - 50%	C = Coarse	R = Rounded			
NA = Not Applicable		some = 20 - 35%	M = Medium	A = Angular			
		little = 10 - 20%	F = Fine	SR = Subrounded			
		trace = 1 - 10%	VF = Very Fine	SA = Subangular			
<b>BORING:</b> SB - 02							

 <b>LaBella</b> Powered by partnership. Soil Boring Log 4 British American Boulevard, Latham, NY			<b>PROJECT</b> Phase II Environmental Site Assessment 5149 County Route 113 Town of Greenwich, Washington County, New York			<b>BORING:</b> SB - 03 <b>SHEET</b> 1 of _1 <b>JOB:</b> 2233938 <b>CHKD BY:</b> BF <b>DATE:</b> 10/24/2023							
CONTRACTOR: Core Down Drilling			BORING LOCATION: South/Downgradient of Anomaly Adjacent "Barn" Building			TIME: 11:00 TO 11:30							
DRILLER:	Joe Bellucci		GROUND SURFACE ELEVATION NA			DATUM: NA							
LABELLA REPRESENTATIVE:	Branson Fields		START DATE:	10/24/23		WEATHER: M.Cloudy, ~40deg.F							
TYPE OF DRILL RIG: Geoprobe 54DT			DRIVE SAMPLER TYPE: Macrocore										
AUGER SIZE AND TYPE: NA			INSIDE DIAMETER: 2-inch										
OVERBURDEN SAMPLING METHOD: Direct Push			OTHER: 4-feet core intervals										
DEPTH (FEET BGS)	SAMPLE		VISUAL CLASSIFICATION			PID FIELD SCREEN (PPM)	REMARKS						
	SAMPLE RECOVERY (INCHES)	SAMPLE NO. AND DEPTH	STRATA CHANGE (FEET BGS)										
0	26/48	No. 01 0 to 4-ft. bgs	Topsoil	6" Topsoil with vegetation			Soil Sample: SB-03(14-15ft) 11:45						
1			SM	20" Very loose, light brown, silty sand (vf-c) with gravel (A), dry, non-plastic, NOSOI.									
2													
3			SM	21" Same as above, moist, NOSOI.									
4	38/48	No. 02 4 to 8-ft. bgs											
5				17" Very loose, light brown, silty sand (f-c) with gravel (A); moist, non-plastic, NOSOI.									
6													
7													
8	38/48	No. 03 8 to 12 ft. bgs.	SM	8" Very loose, dark brown, silty sand (f-c), moist, little gravel (A) non-plastic, NOSOI.			<1						
9				30" Very loose, dark brown/red, silty gravel (A) with sand (f-c), moist, non-plastic, NOSOI.									
10			GM										
11													
12	38/48	No. 04 12 to 16-ft. bgs	GM	38" Very loose, red/orange, silty gravel (A/SR), moist, non-plastic, little sand (f-c), NOSOI			<1						
13													
14													
15													
16				End of Boring at 16-ft. bgs.									
WATER LEVEL DATA			DEPTH (FT)	NOTES:									
DATE	TIME	ELAPSED TIME	BOTTOM OF CASING	BOTTOM OF BORING	GROUNDWATER ENCOUNTERED								
					na								
GENERAL NOTES													
1) STRATIFICATION LINES REPRESENT APPROXIMATE BOUNDARY BETWEEN SOIL TYPES, TRANSITIONS MAY BE GRADUAL.													
2) WATER LEVEL READINGS HAVE BEEN MADE AT TIMES AND UNDER CONDITIONS STATED, FLUCTUATIONS OF GROUNDWATER													
BGS = Below Ground Surface			and = 35 - 50%	C = Coarse	R = Rounded								
NA = Not Applicable			some = 20 - 35%	M = Medium	A = Angular								
			little = 10 - 20%	F = Fine	SR = Subrounded								
			trace = 1 - 10%	VF = Very Fine	SA = Subangular								
								<b>BORING:</b> SB - 03					



## APPENDIX 2

Laboratory Analytical Report



Thursday, November 09, 2023

Attn: Branson Fields  
Labella Assosiates  
4 British American Blvd  
Latham, NY 12110

Project ID: T.GREENWICH-5149 CR 113  
SDG ID: GCP33746  
Sample ID#s: CP33746 - CP33748

This laboratory is in compliance with the NELAC requirements of procedures used except where indicated.

This report contains results for the parameters tested, under the sampling conditions described on the Chain Of Custody, as received by the laboratory. This report is incomplete unless all pages indicated in the pagination at the bottom of the page are included.

All soils, solids and sludges are reported on a dry weight basis unless otherwise noted in the sample comments.

A scanned version of the COC form accompanies the analytical report and is an exact duplicate of the original.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.

Sincerely yours,

A handwritten signature in black ink, appearing to read "Phyllis Shiller".

Phyllis Shiller

Laboratory Director

NELAC - #NY11301  
CT Lab Registration #PH-0618  
MA Lab Registration #M-CT007  
ME Lab Registration #CT-007  
NH Lab Registration #213693-A,B

NJ Lab Registration #CT-003  
NY Lab Registration #11301  
PA Lab Registration #68-03530  
RI Lab Registration #63  
VT Lab Registration #VT11301



Environmental Laboratories, Inc.  
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102 Fax (860) 645-0823



## SDG Comments

November 09, 2023

SDG I.D.: GCP33746

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Please be advised that the NY 375 soil criteria for chromium are based on hexavalent chromium and trivalent chromium.



Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102 Fax (860) 645-0823

## Sample Id Cross Reference

November 09, 2023

SDG I.D.: GCP33746

Project ID: T.GREENWICH-5149 CR 113

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Client Id	Lab Id	Matrix
SB-01 (6-7FT)	CP33746	SOIL
SB-02 (7-8FT)	CP33747	SOIL
SB-03 (14-15FT)	CP33748	SOIL



## Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102 Fax (860) 645-0823



# Analysis Report

November 09, 2023

FOR: Attn: Branson Fields  
Labella Assosiates  
4 British American Blvd  
Latham, NY 12110

### Sample Information

Matrix: SOIL  
Location Code: LABELLA-LATHAM  
Rush Request: Standard  
P.O.#: 2233938

### Custody Information

Collected by: BF  
Received by: SR1  
Analyzed by: see "By" below

Date

10/24/23 9:55  
10/25/23 17:06

Time

Project ID: T.GREENWICH-5149 CR 113  
Client ID: SB-01 (6-7FT)

### Laboratory Data

SDG ID: GCP33746

Phoenix ID: CP33746

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.44	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Aluminum	17100	67	mg/Kg	10	10/31/23	IE	SW6010D
Arsenic	4.08	0.89	mg/Kg	1	10/31/23	IE	SW6010D
Barium	122	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Beryllium	0.82	0.36	mg/Kg	1	10/31/23	IE	SW6010D
Calcium	4900	6.7	mg/Kg	1	10/31/23	IE	SW6010D
Cadmium	0.91	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Cobalt	11.4	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Chromium	18.3	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Copper	15.1	0.9	mg/kg	1	10/31/23	IE	SW6010D
Iron	33600	67	mg/Kg	10	10/31/23	IE	SW6010D
Mercury	< 0.03	0.03	mg/Kg	2	10/26/23	AL1	SW7471B
Potassium	1390	67	mg/Kg	10	10/31/23	IE	SW6010D
Magnesium	4260	6.7	mg/Kg	1	10/31/23	IE	SW6010D
Manganese	340	4.4	mg/Kg	10	10/31/23	IE	SW6010D
Sodium	263	6.7	mg/Kg	1	10/31/23	IE	SW6010D
Nickel	17.2	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Lead	9.00	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Antimony	< 4.4	4.4	mg/Kg	1	10/31/23	IE	SW6010D
Selenium	< 1.8	1.8	mg/Kg	1	10/31/23	IE	SW6010D
Thallium	< 4.0	4.0	mg/Kg	1	10/31/23	IE	SW6010D
Vanadium	36.6	0.44	mg/Kg	1	10/31/23	IE	SW6010D
Zinc	73.5	0.9	mg/Kg	1	10/31/23	IE	SW6010D
Percent Solid	77		%		10/25/23	CV	SW846-%Solid
Mercury Digestion	Completed				10/26/23	AL/AL	SW7471B
Soil Extraction for SVOA	Completed				10/30/23	B/A	SW3546
Total Metals Digest	Completed				10/26/23	B/AG	SW3050B

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
<b>Volatiles (TCL)</b>							
1,1,1-Trichloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,1,2,2-Tetrachloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,1,2-Trichloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,1-Dichloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,1-Dichloroethene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2,3-Trichlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2,4-Trichlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dibromo-3-chloropropane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dibromoethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichloropropane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,3-Dichlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
1,4-Dichlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
2-Hexanone	ND	0.028	mg/Kg	1	10/26/23	JLI	SW8260D
4-Methyl-2-pentanone	ND	0.028	mg/Kg	1	10/26/23	JLI	SW8260D
Acetone	ND	0.05	mg/Kg	1	10/26/23	JLI	SW8260D
Benzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Bromochloromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Bromodichloromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Bromoform	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Bromomethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Carbon Disulfide	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Carbon tetrachloride	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Chlorobenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Chloroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Chloroform	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Chloromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
cis-1,2-Dichloroethene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
cis-1,3-Dichloropropene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Cyclohexane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Dibromochloromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Dichlorodifluoromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Ethylbenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Isopropylbenzene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
m&p-Xylene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Methyl ethyl ketone	ND	0.034	mg/Kg	1	10/26/23	JLI	SW8260D
Methyl t-butyl ether (MTBE)	ND	0.011	mg/Kg	1	10/26/23	JLI	SW8260D
Methylacetate	ND	0.0045	mg/Kg	1	10/26/23	JLI	SW8260D
Methylcyclohexane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Methylene chloride	ND	0.028	mg/Kg	1	10/26/23	JLI	SW8260D
o-Xylene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Styrene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Tetrachloroethene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Toluene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Total Xylenes	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
trans-1,2-Dichloroethene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
trans-1,3-Dichloropropene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Trichloroethene	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Trichlorofluoromethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Trichlorotrifluoroethane	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
Vinyl chloride	ND	0.0056	mg/Kg	1	10/26/23	JLI	SW8260D
<b><u>QA/QC Surrogates</u></b>							
% 1,2-dichlorobenzene-d4	96		%	1	10/26/23	JLI	70 - 130 %
% Bromofluorobenzene	94		%	1	10/26/23	JLI	70 - 130 %
% Dibromofluoromethane	96		%	1	10/26/23	JLI	70 - 130 %
% Toluene-d8	91		%	1	10/26/23	JLI	70 - 130 %
<b><u>1,4-dioxane</u></b>							
1,4-dioxane	ND	0.084	mg/Kg	1	10/26/23	JLI	SW8260D
<b><u>Semivolatiles</u></b>							
1,1-Biphenyl	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
1,2,4,5-Tetrachlorobenzene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,2'-Oxybis(1-Chloropropane)	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,3,4,6-tetrachlorophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,4,5-Trichlorophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,4,6-Trichlorophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dichlorophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dimethylphenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dinitrophenol	ND	0.67	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dinitrotoluene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2,6-Dinitrotoluene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2-Chloronaphthalene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2-Chlorophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2-Methylnaphthalene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2-Methylphenol (o-cresol)	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
2-Nitroaniline	ND	0.67	mg/Kg	1	10/30/23	KCA	SW8270D
2-Nitrophenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
3,3'-Dichlorobenzidine	ND	0.5	mg/Kg	1	10/30/23	KCA	SW8270D
3-Nitroaniline	ND	0.67	mg/Kg	1	10/30/23	KCA	SW8270D
4,6-Dinitro-2-methylphenol	ND	1.2	mg/Kg	1	10/30/23	KCA	SW8270D
4-Bromophenyl phenyl ether	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chloro-3-methylphenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chloroaniline	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chlorophenyl phenyl ether	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
4-Nitroaniline	ND	0.67	mg/Kg	1	10/30/23	KCA	SW8270D
4-Nitrophenol	ND	1.2	mg/Kg	1	10/30/23	KCA	SW8270D
Acenaphthene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Acenaphthylene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Acetophenone	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Anthracene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Atrazine	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benz(a)anthracene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benzaldehyde	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(a)pyrene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(b)fluoranthene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Benzo(ghi)perylene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(k)fluoranthene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Benzyl butyl phthalate	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-chloroethoxy)methane	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-chloroethyl)ether	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-ethylhexyl)phthalate	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Caprolactam	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Carbazole	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
Chrysene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Dibenz(a,h)anthracene	ND	0.21	mg/Kg	1	10/30/23	KCA	SW8270D
Dibenzofuran	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Diethyl phthalate	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Dimethylphthalate	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Di-n-butylphthalate	ND	0.84	mg/Kg	1	10/30/23	KCA	SW8270D
Di-n-octylphthalate	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Fluoranthene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Fluorene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorobenzene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorobutadiene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorocyclopentadiene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachloroethane	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Indeno(1,2,3-cd)pyrene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Isophorone	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Naphthalene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Nitrobenzene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodimethylamine	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodi-n-propylamine	ND	0.21	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodiphenylamine	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
Pentachlorophenol	ND	0.42	mg/Kg	1	10/30/23	KCA	SW8270D
Phenanthrene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Phenol	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
Pyrene	ND	0.29	mg/Kg	1	10/30/23	KCA	SW8270D
<b><u>QA/QC Surrogates</u></b>							
% 2,4,6-Tribromophenol	71		%	1	10/30/23	KCA	30 - 130 %
% 2-Fluorobiphenyl	62		%	1	10/30/23	KCA	30 - 130 %
% 2-Fluorophenol	59		%	1	10/30/23	KCA	30 - 130 %
% Nitrobenzene-d5	56		%	1	10/30/23	KCA	30 - 130 %
% Phenol-d5	60		%	1	10/30/23	KCA	30 - 130 %
% Terphenyl-d14	67		%	1	10/30/23	KCA	30 - 130 %

SVOA Library Search Top 15

Completed

11/08/23

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Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
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1 = This parameter is not certified by the primary accrediting authority (NY NELAC) for this matrix. NY NELAC does not offer certification for all parameters at this time.

RL/PQL=Reporting/Practical Quantitation Level (Equivalent to NELAC LOQ, Limit of Quantitation) ND=Not Detected at RL/PQL  
BRL=Below Reporting Level L=Biased Low

QA/QC Surrogates: Surrogates are compounds (preceded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

### **Comments:**

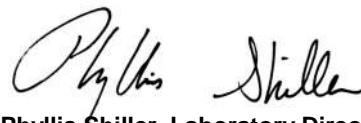
Please be advised that the NY 375 soil criteria for chromium are based on hexavalent chromium and trivalent chromium.

#### Volatile Comment:

To achieve client's objectives, where the lowest calibration standard or LOD justifies lowering the RL/PQL, the RL/PQL of some compounds have been lowered to meet criteria.

All soils, solids and sludges are reported on a dry weight basis unless otherwise noted in the sample comments.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.



Phyllis Shiller, Laboratory Director

November 09, 2023

Reviewed and Released by: Phyllis Shiller, Laboratory Director



## Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102 Fax (860) 645-0823



# Analysis Report

November 09, 2023

FOR: Attn: Branson Fields  
Labella Assosiates  
4 British American Blvd  
Latham, NY 12110

### Sample Information

Matrix: SOIL  
Location Code: LABELLA-LATHAM  
Rush Request: Standard  
P.O.#: 2233938

### Custody Information

Collected by: BF  
Received by: SR1  
Analyzed by: see "By" below

Date

Time

SDG ID: GCP33746

Phoenix ID: CP33747

Project ID: T.GREENWICH-5149 CR 113  
Client ID: SB-02 (7-8FT)

### Laboratory Data

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.39	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Aluminum	32600	5800	mg/Kg	1000	11/07/23	TH	SW6010D
Arsenic	4.93	0.78	mg/Kg	1	10/31/23	IE	SW6010D
Barium	90.5	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Beryllium	0.97	0.31	mg/Kg	1	10/31/23	IE	SW6010D
Calcium	1380	5.8	mg/Kg	1	10/31/23	IE	SW6010D
Cadmium	1.99	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Cobalt	15.0	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Chromium	10.8	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Copper	36.6	0.8	mg/kg	1	10/31/23	IE	SW6010D
Iron	69100	58	mg/Kg	10	10/31/23	IE	SW6010D
Mercury	< 0.03	0.03	mg/Kg	2	10/26/23	AL1	SW7471B
Potassium	620	5.8	mg/Kg	1	10/31/23	IE	SW6010D
Magnesium	11200	58	mg/Kg	10	10/31/23	IE	SW6010D
Manganese	1140	3.9	mg/Kg	10	10/31/23	IE	SW6010D
Sodium	28.2	5.8	mg/Kg	1	10/31/23	IE	SW6010D
Nickel	55.9	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Lead	11.6	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Antimony	< 3.9	3.9	mg/Kg	1	10/31/23	IE	SW6010D
Selenium	< 1.6	1.6	mg/Kg	1	10/31/23	IE	SW6010D
Thallium	< 3.5	3.5	mg/Kg	1	10/31/23	IE	SW6010D
Vanadium	23.2	0.39	mg/Kg	1	10/31/23	IE	SW6010D
Zinc	174	7.8	mg/Kg	10	10/31/23	IE	SW6010D
Percent Solid	87		%		10/25/23	CV	SW846-%Solid
Mercury Digestion	Completed				10/26/23	AL/AL	SW7471B
Soil Extraction for SVOA	Completed				10/30/23	B/A	SW3546
Total Metals Digest	Completed				10/26/23	B/AG	SW3050B

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
<b>Volatiles (TCL)</b>							
1,1,1-Trichloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,1,2,2-Tetrachloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,1,2-Trichloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,1-Dichloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,1-Dichloroethene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2,3-Trichlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2,4-Trichlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dibromo-3-chloropropane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dibromoethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,2-Dichloropropane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,3-Dichlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
1,4-Dichlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
2-Hexanone	ND	0.017	mg/Kg	1	10/26/23	JLI	SW8260D
4-Methyl-2-pentanone	ND	0.017	mg/Kg	1	10/26/23	JLI	SW8260D
Acetone	ND	0.034	mg/Kg	1	10/26/23	JLI	SW8260D
Benzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Bromochloromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Bromodichloromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Bromoform	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Bromomethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Carbon Disulfide	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Carbon tetrachloride	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Chlorobenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Chloroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Chloroform	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Chloromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
cis-1,2-Dichloroethene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
cis-1,3-Dichloropropene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Cyclohexane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Dibromochloromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Dichlorodifluoromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Ethylbenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Isopropylbenzene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
m&p-Xylene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Methyl ethyl ketone	ND	0.021	mg/Kg	1	10/26/23	JLI	SW8260D
Methyl t-butyl ether (MTBE)	ND	0.0069	mg/Kg	1	10/26/23	JLI	SW8260D
Methylacetate	ND	0.0028	mg/Kg	1	10/26/23	JLI	SW8260D
Methylcyclohexane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Methylene chloride	ND	0.017	mg/Kg	1	10/26/23	JLI	SW8260D
o-Xylene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Styrene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Tetrachloroethene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Toluene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Total Xylenes	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
trans-1,2-Dichloroethene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
trans-1,3-Dichloropropene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Trichloroethene	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Trichlorofluoromethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Trichlorotrifluoroethane	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
Vinyl chloride	ND	0.0034	mg/Kg	1	10/26/23	JLI	SW8260D
<b><u>QA/QC Surrogates</u></b>							
% 1,2-dichlorobenzene-d4	97		%	1	10/26/23	JLI	70 - 130 %
% Bromofluorobenzene	94		%	1	10/26/23	JLI	70 - 130 %
% Dibromofluoromethane	96		%	1	10/26/23	JLI	70 - 130 %
% Toluene-d8	91		%	1	10/26/23	JLI	70 - 130 %
<b><u>1,4-dioxane</u></b>							
1,4-dioxane	ND	0.052	mg/Kg	1	10/26/23	JLI	SW8260D
<b><u>Semivolatiles</u></b>							
1,1-Biphenyl	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
1,2,4,5-Tetrachlorobenzene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,2'-Oxybis(1-Chloropropane)	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,3,4,6-tetrachlorophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,4,5-Trichlorophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,4,6-Trichlorophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dichlorophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dimethylphenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dinitrophenol	ND	0.6	mg/Kg	1	10/30/23	KCA	SW8270D
2,4-Dinitrotoluene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2,6-Dinitrotoluene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2-Chloronaphthalene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2-Chlorophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2-Methylnaphthalene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2-Methylphenol (o-cresol)	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
2-Nitroaniline	ND	0.6	mg/Kg	1	10/30/23	KCA	SW8270D
2-Nitrophenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
3,3'-Dichlorobenzidine	ND	0.45	mg/Kg	1	10/30/23	KCA	SW8270D
3-Nitroaniline	ND	0.6	mg/Kg	1	10/30/23	KCA	SW8270D
4,6-Dinitro-2-methylphenol	ND	1.1	mg/Kg	1	10/30/23	KCA	SW8270D
4-Bromophenyl phenyl ether	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chloro-3-methylphenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chloroaniline	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
4-Chlorophenyl phenyl ether	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
4-Nitroaniline	ND	0.6	mg/Kg	1	10/30/23	KCA	SW8270D
4-Nitrophenol	ND	1.1	mg/Kg	1	10/30/23	KCA	SW8270D
Acenaphthene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Acenaphthylene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Acetophenone	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Anthracene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Atrazine	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benz(a)anthracene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benzaldehyde	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(a)pyrene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(b)fluoranthene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Benzo(ghi)perylene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benzo(k)fluoranthene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Benzyl butyl phthalate	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-chloroethoxy)methane	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-chloroethyl)ether	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
Bis(2-ethylhexyl)phthalate	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Caprolactam	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Carbazole	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
Chrysene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Dibenz(a,h)anthracene	ND	0.19	mg/Kg	1	10/30/23	KCA	SW8270D
Dibenzofuran	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Diethyl phthalate	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Dimethylphthalate	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Di-n-butylphthalate	ND	0.75	mg/Kg	1	10/30/23	KCA	SW8270D
Di-n-octylphthalate	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Fluoranthene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Fluorene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorobenzene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorobutadiene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachlorocyclopentadiene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Hexachloroethane	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Indeno(1,2,3-cd)pyrene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Isophorone	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Naphthalene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Nitrobenzene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodimethylamine	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodi-n-propylamine	ND	0.19	mg/Kg	1	10/30/23	KCA	SW8270D
N-Nitrosodiphenylamine	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
Pentachlorophenol	ND	0.37	mg/Kg	1	10/30/23	KCA	SW8270D
Phenanthrene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Phenol	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
Pyrene	ND	0.26	mg/Kg	1	10/30/23	KCA	SW8270D
<b><u>QA/QC Surrogates</u></b>							
% 2,4,6-Tribromophenol	71		%	1	10/30/23	KCA	30 - 130 %
% 2-Fluorobiphenyl	61		%	1	10/30/23	KCA	30 - 130 %
% 2-Fluorophenol	58		%	1	10/30/23	KCA	30 - 130 %
% Nitrobenzene-d5	55		%	1	10/30/23	KCA	30 - 130 %
% Phenol-d5	60		%	1	10/30/23	KCA	30 - 130 %
% Terphenyl-d14	68		%	1	10/30/23	KCA	30 - 130 %

SVOA Library Search Top 15

Completed

11/08/23

E

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
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1 = This parameter is not certified by the primary accrediting authority (NY NELAC) for this matrix. NY NELAC does not offer certification for all parameters at this time.

RL/PQL=Reporting/Practical Quantitation Level (Equivalent to NELAC LOQ, Limit of Quantitation) ND=Not Detected at RL/PQL  
BRL=Below Reporting Level L=Biased Low

QA/QC Surrogates: Surrogates are compounds (preceded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

### **Comments:**

Please be advised that the NY 375 soil criteria for chromium are based on hexavalent chromium and trivalent chromium.

All soils, solids and sludges are reported on a dry weight basis unless otherwise noted in the sample comments.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.



Phyllis Shiller, Laboratory Director

November 09, 2023

Reviewed and Released by: Phyllis Shiller, Laboratory Director



## Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102 Fax (860) 645-0823



# Analysis Report

November 09, 2023

FOR: Attn: Branson Fields  
Labella Assosiates  
4 British American Blvd  
Latham, NY 12110

### Sample Information

Matrix: SOIL  
Location Code: LABELLA-LATHAM  
Rush Request: Standard  
P.O.#: 2233938

### Custody Information

Collected by: BF  
Received by: SR1  
Analyzed by: see "By" below

Date

Time

10/24/23 11:45

10/25/23 17:06

SDG ID: GCP33746

Phoenix ID: CP33748

Project ID: T.GREENWICH-5149 CR 113  
Client ID: SB-03 (14-15FT)

### Laboratory Data

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
Percent Solid	89		%		10/25/23	CV	SW846-%Solid
Soil Extraction for SVOA PAH	Completed				10/27/23	H/JDW	SW3546
<b>Volatiles- STARS/CP-51</b>							
1,2,4-Trimethylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
1,3,5-Trimethylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
Benzene	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
Ethylbenzene	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
Isopropylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
m&p-Xylene	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
Methyl t-Butyl Ether (MTBE)	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
Naphthalene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
n-Butylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
n-Propylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
o-Xylene	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
p-Isopropyltoluene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
sec-Butylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
tert-Butylbenzene	ND	0.00093	mg/Kg	1	10/26/23	JLI	SW8260D
Toluene	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
Total Xylenes	ND	0.0019	mg/Kg	1	10/26/23	JLI	SW8260D
<b>QA/QC Surrogates</b>							
% 1,2-Dichlorobenzene-d4	100		%	1	10/26/23	JLI	70 - 130 %
% Bromofluorobenzene	96		%	1	10/26/23	JLI	70 - 130 %
% Dibromofluoromethane	98		%	1	10/26/23	JLI	70 - 130 %
% Toluene-d8	98		%	1	10/26/23	JLI	70 - 130 %

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	By	Reference
<b>Semivolatiles-STARS/CP-51</b>							
Acenaphthene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Acenaphthylene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Anthracene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Benz(a)anthracene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Benzo(a)pyrene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Benzo(b)fluoranthene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Benzo(ghi)perylene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Benzo(k)fluoranthene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Chrysene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Dibenz(a,h)anthracene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Fluoranthene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Fluorene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Indeno(1,2,3-cd)pyrene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Naphthalene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Phenanthrene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
Pyrene	ND	0.26	mg/Kg	1	10/29/23	KCA	SW8270D
<b>QA/QC Surrogates</b>							
% 2-Fluorobiphenyl	57		%	1	10/29/23	KCA	30 - 130 %
% Nitrobenzene-d5	57		%	1	10/29/23	KCA	30 - 130 %
% Terphenyl-d14	61		%	1	10/29/23	KCA	30 - 130 %

RL/PQL=Reporting/Practical Quantitation Level (Equivalent to NELAC LOQ, Limit of Quantitation) ND=Not Detected at RL/PQL  
 BRL=Below Reporting Level L=Biased Low

QA/QC Surrogates: Surrogates are compounds (preceded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

### Comments:

All soils, solids and sludges are reported on a dry weight basis unless otherwise noted in the sample comments.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.

Phyllis Shiller, Laboratory Director

November 09, 2023

Reviewed and Released by: Phyllis Shiller, Laboratory Director

**1F**  
**SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET**  
**TENTATIVELY IDENTIFIED COMPOUNDS**

## CLIENT ID

SB-01 (6-7FT)

Lab Name: Phoenix Environmental Labs

Client: LABELLA-LATHAM

Lab Code: Phoenix Case No.:

SAS No.: \_\_\_\_\_ SDG No.: GCP33746

## Matrix:(soil/water)            SOIL

Lab Sample ID: CP33746

Sample wt/vol: 15.46 (g/mL) g

Lab File ID: 1030\_13.D

Level: (low/med)  Low

Date Received: 10/25/23

% Moisture: not dec. 23 decanted:(Y/N) NA

Date Extracted: 10/30/23

GPC Cleanup (Y/N): N pH: NA

Date Analyzed: 10/30/2023

Conc. Extract Volume: 1000 (uL)

Dilution Factor 1

Conc. Extract Volume: 1000 (uL) Dilution Factor 1

Injection Volume: 1 (uL)

Injection Volume: 1 (uL)

**CONCENTRATION UNITS:**

(ug/L or ug/KG)      ug/Kg

## FORM I SEMIVOYA-TIC

- A - Indicates that the tentatively identified compound is a suspected aldol condensation product.  
Aldol condensation products are produced during the extraction process.
  - C - Indicates that the tentatively identified compound is a suspected prep artifact produced during extraction process.
  - Q - For TICS, this compound was quantitated using a calibration curve. This compound is part of the instrument method, but not part of the client target list.

**1F**  
**SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET**  
**TENTATIVELY IDENTIFIED COMPOUNDS**

## CLIENT ID

SB-02 (7-8FT)

Lab Name: Phoenix Environmental Labs

Client: LABELLA-LATHAM

Lab Code: Phoenix Case No.:

SAS No.: \_\_\_\_\_ SDG No.: GCP33746

SDG No.: GCP33746

## Matrix:(soil/water)            SOIL

Lab Sample ID: CP33747

Sample wt/vol: 15.38 (g/mL) g

Lab File ID: 1030\_14.D

Level: (low/med)  Low

Date Received: 10/25/23

% Moisture: not dec. 13 decanted:(Y/N) NA

Date Extracted: 10/30/23

GPC Cleanup (Y/N): N pH: NA

Date Analyzed: 10/30/2023

Conc. Extract Volume: 1000 (uL)

Dilution Factor 1

Conc. Extract Volume: 1000 (uL) Dilution Factor 1

Injection Volume: 1 (uL)

Injection Volume: 1 ( $\mu\text{L}$ )

**CONCENTRATION UNITS:**

Number TICs found: 4

**CONCENTRATION UNITS:**

Number TICs found: 4 (ug/L or ug/KG) ug/Kg

FORM I SEMIVOYA-TIC

- A - Indicates that the tentatively identified compound is a suspected aldol condensation product.  
Aldol condensation products are produced during the extraction process.
  - C - Indicates that the tentatively identified compound is a suspected prep artifact produced during extraction process.
  - Q - For TICS, this compound was quantitated using a calibration curve. This compound is part of the instrument method, but not part of the client target list.



**Environmental Laboratories, Inc.**

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045

Tel. (860) 645-1102



## QA/QC Report

November 09, 2023

### QA/QC Data

SDG I.D.: GCP33746

Parameter	Blank	Blk RL	Sample Result	Dup Result	Dup RPD	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
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QA/QC Batch 703610 (mg/kg), QC Sample No: CP33802 2X (CP33746, CP33747)

Mercury - Soil	BRL	0.03	<0.03	<0.03	NC	123	122	0.8	91.4	94.9	3.8	70 - 130	30
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Comment:

Additional Mercury criteria: LCS acceptance range for waters is 80-120% and for soils is 70-130%. MS acceptance range is 75-125%.

QA/QC Batch 703536 (mg/kg), QC Sample No: CP33454 (CP33746, CP33747)

### ICP Metals - Soil

Aluminum	BRL	5.0	11900	11200	6.10	105	102	2.9	NC		75 - 125	35
Antimony	BRL	3.3	5.8	5.0	NC	111	112	0.9	86.0		75 - 125	35
Arsenic	BRL	0.67	13.0	11.1	15.8	100	97.1	2.9	98.1		75 - 125	35
Barium	BRL	0.33	141	130	8.10	102	99.7	2.3	94.2		75 - 125	35
Beryllium	BRL	0.27	0.52	0.47	NC	98.3	93.6	4.9	97.0		75 - 125	35
Cadmium	BRL	0.33	5.33	3.84	32.5	105	98.5	6.4	97.9		75 - 125	35
Calcium	BRL	5.0	6710	4780	33.6	101	96.5	4.6	NC		75 - 125	35
Chromium	BRL	0.33	69.4	78.9	12.8	102	99.5	2.5	92.1		75 - 125	35
Cobalt	BRL	0.33	11.4	9.69	16.2	101	98.8	2.2	98.2		75 - 125	35
Copper	BRL	0.67	134	133	0.70	95.7	94.8	0.9	92.9		75 - 125	35
Iron	BRL	5.0	53200	50600	5.00	105	105	0.0	NC		75 - 125	35
Lead	BRL	0.33	663	192	110	102	100	2.0	105		75 - 125	35
Magnesium	BRL	5.0	5460	4660	15.8	108	104	3.8	NC		75 - 125	35
Manganese	BRL	0.33	457	444	2.90	112	99.1	12.2	82.7		75 - 125	35
Nickel	BRL	0.33	46.7	205	126	103	99.6	3.4	94.3		75 - 125	35
Potassium	BRL	5.0	2930	2800	4.50	90.4	91.6	1.3	123		75 - 125	35
Selenium	BRL	1.3	<1.5	<1.4	NC	102	98.7	3.3	94.4		75 - 125	35
Silver	BRL	0.33	<3.9	1.00	NC	96.8	95.3	1.6	88.1		75 - 125	35
Sodium	BRL	5.0	278	300	7.60	83.3	84.8	1.8	113		75 - 125	35
Thallium	BRL	3.0	<3.4	<3.1	NC	101	96.7	4.4	97.0		75 - 125	35
Vanadium	BRL	0.33	38.7	35.7	8.10	101	99.2	1.8	100		75 - 125	35
Zinc	BRL	0.67	745	624	17.7	96.0	94.1	2.0	NC		75 - 125	35

Comment:

Additional Criteria: LCS acceptance range is 80-120% MS acceptance range 75-125%.

r = This parameter is outside laboratory RPD specified recovery limits.



Environmental Laboratories, Inc.  
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
Tel. (860) 645-1102

## QA/QC Report

November 09, 2023

### QA/QC Data

SDG I.D.: GCP33746

Parameter	Blank	Blk	RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
QA/QC Batch 703956 (mg/Kg), QC Sample No: CP33748 (CP33748)											
<u>Polynuclear Aromatic HC - Soil</u>											
Acenaphthene	ND	0.23		68	66	3.0	56	61	8.5	30 - 130	30
Acenaphthylene	ND	0.23		66	62	6.3	55	59	7.0	40 - 140	30
Anthracene	ND	0.23		70	67	4.4	60	63	4.9	40 - 140	30
Benz(a)anthracene	ND	0.23		67	66	1.5	59	63	6.6	40 - 140	30
Benzo(a)pyrene	ND	0.23		74	69	7.0	63	67	6.2	40 - 140	30
Benzo(b)fluoranthene	ND	0.23		68	66	3.0	59	62	5.0	40 - 140	30
Benzo(ghi)perylene	ND	0.23		76	70	8.2	64	67	4.6	40 - 140	30
Benzo(k)fluoranthene	ND	0.23		70	70	0.0	63	63	0.0	40 - 140	30
Chrysene	ND	0.23		70	69	1.4	59	65	9.7	40 - 140	30
Dibenz(a,h)anthracene	ND	0.23		74	69	7.0	62	65	4.7	40 - 140	30
Fluoranthene	ND	0.23		73	79	7.9	70	69	1.4	40 - 140	30
Fluorene	ND	0.23		67	72	7.2	63	67	6.2	40 - 140	30
Indeno(1,2,3-cd)pyrene	ND	0.23		73	69	5.6	62	66	6.3	40 - 140	30
Naphthalene	ND	0.23		63	62	1.6	55	57	3.6	40 - 140	30
Phenanthrene	ND	0.23		68	66	3.0	58	62	6.7	40 - 140	30
Pyrene	ND	0.23		73	80	9.2	71	69	2.9	30 - 130	30
% 2-Fluorobiphenyl	68	%		70	57	20.5	51	55	7.5	30 - 130	30
% Nitrobenzene-d5	71	%		73	61	17.9	59	60	1.7	30 - 130	30
% Terphenyl-d14	73	%		67	76	12.6	65	66	1.5	30 - 130	30
QA/QC Batch 704104 (mg/Kg), QC Sample No: CP35727 (CP33746, CP33747)											
<u>Semivolatiles - Soil</u>											
1,1-Biphenyl	ND	0.23		60	61	1.7	60	57	5.1	40 - 140	30
1,2,4,5-Tetrachlorobenzene	ND	0.23		62	63	1.6	63	58	8.3	40 - 140	30
2,2'-Oxybis(1-Chloropropane)	ND	0.23		42	45	6.9	45	41	9.3	40 - 140	30
2,3,4,6-tetrachlorophenol	ND	0.23		76	75	1.3	79	68	15.0	30 - 130	30
2,4,5-Trichlorophenol	ND	0.23		76	76	0.0	77	70	9.5	40 - 140	30
2,4,6-Trichlorophenol	ND	0.13		76	76	0.0	78	71	9.4	30 - 130	30
2,4-Dichlorophenol	ND	0.13		71	72	1.4	76	67	12.6	30 - 130	30
2,4-Dimethylphenol	ND	0.23		70	70	0.0	76	67	12.6	30 - 130	30
2,4-Dinitrophenol	ND	0.23		80	84	4.9	84	58	36.6	30 - 130	30
2,4-Dinitrotoluene	ND	0.13		76	76	0.0	80	71	11.9	30 - 130	30
2,6-Dinitrotoluene	ND	0.13		75	75	0.0	78	70	10.8	40 - 140	30
2-Chloronaphthalene	ND	0.23		62	63	1.6	62	59	5.0	40 - 140	30
2-Chlorophenol	ND	0.23		61	67	9.4	69	61	12.3	30 - 130	30
2-Methylnaphthalene	ND	0.23		64	65	1.6	66	61	7.9	40 - 140	30
2-Methylphenol (o-cresol)	ND	0.23		58	63	8.3	68	58	15.9	40 - 140	30
2-Nitroaniline	ND	0.33		83	87	4.7	98	84	15.4	40 - 140	30
2-Nitrophenol	ND	0.23		70	73	4.2	75	69	8.3	40 - 140	30
3&4-Methylphenol (m&p-cresol)	ND	0.23		64	69	7.5	76	63	18.7	30 - 130	30
3,3'-Dichlorobenzidine	ND	0.13		55	98	56.2	99	87	12.9	40 - 140	30
3-Nitroaniline	ND	0.33		67	86	24.8	96	81	16.9	40 - 140	30

QA/QC Data

SDG I.D.: GCP33746

Parameter	Blank	Blk	RL	LCS	LCSD	LCS	MS	MSD	MS	%	%
				%	%	RPD	%	MSD %	MS RPD	Rec Limits	RPD Limits
4,6-Dinitro-2-methylphenol	ND	0.23		84	86	2.4	89	70	23.9	30 - 130	30
4-Bromophenyl phenyl ether	ND	0.23		71	71	0.0	71	67	5.8	40 - 140	30
4-Chloro-3-methylphenol	ND	0.23		73	74	1.4	81	69	16.0	30 - 130	30
4-Chloroaniline	ND	0.23		50	58	14.8	70	57	20.5	40 - 140	30
4-Chlorophenyl phenyl ether	ND	0.23		68	68	0.0	69	63	9.1	40 - 140	30
4-Nitroaniline	ND	0.23		71	71	0.0	77	67	13.9	40 - 140	30
4-Nitrophenol	ND	0.23		86	96	11.0	109	89	20.2	30 - 130	30
Acenaphthene	ND	0.23		64	64	0.0	65	59	9.7	30 - 130	30
Acenaphthylene	ND	0.13		61	62	1.6	63	59	6.6	40 - 140	30
Acetophenone	ND	0.23		54	58	7.1	59	52	12.6	40 - 140	30
Anthracene	ND	0.23		68	68	0.0	70	65	7.4	40 - 140	30
Atrazine	ND	0.13		58	69	17.3	73	68	7.1	40 - 140	30
Benz(a)anthracene	ND	0.23		68	66	3.0	66	61	7.9	40 - 140	30
Benzaldehyde	ND	0.23		63	66	4.7	71	45	44.8	40 - 140	30
Benzo(a)pyrene	ND	0.13		78	74	5.3	72	66	8.7	40 - 140	30
Benzo(b)fluoranthene	ND	0.16		74	66	11.4	65	59	9.7	40 - 140	30
Benzo(ghi)perylene	ND	0.23		78	70	10.8	71	67	5.8	40 - 140	30
Benzo(k)fluoranthene	ND	0.23		74	67	9.9	67	62	7.8	40 - 140	30
Benzyl butyl phthalate	ND	0.23		64	65	1.6	65	60	8.0	40 - 140	30
Bis(2-chloroethoxy)methane	ND	0.23		58	60	3.4	61	55	10.3	40 - 140	30
Bis(2-chloroethyl)ether	ND	0.13		54	53	1.9	57	48	17.1	40 - 140	30
Bis(2-ethylhexyl)phthalate	ND	0.23		74	75	1.3	75	71	5.5	40 - 140	30
Caprolactam	ND	0.23		73	73	0.0	85	67	23.7	40 - 140	30
Carbazole	ND	0.23		69	69	0.0	73	66	10.1	40 - 140	30
Chrysene	ND	0.23		69	67	2.9	66	62	6.3	40 - 140	30
Dibenz(a,h)anthracene	ND	0.13		77	69	11.0	69	65	6.0	40 - 140	30
Dibenzofuran	ND	0.23		64	63	1.6	65	60	8.0	40 - 140	30
Diethyl phthalate	ND	0.23		68	67	1.5	71	62	13.5	40 - 140	30
Dimethylphthalate	ND	0.23		68	67	1.5	70	62	12.1	40 - 140	30
Di-n-butylphthalate	ND	0.67		74	72	2.7	74	67	9.9	40 - 140	30
Di-n-octylphthalate	ND	0.23		66	68	3.0	68	63	7.6	40 - 140	30
Fluoranthene	ND	0.23		72	71	1.4	70	62	12.1	40 - 140	30
Fluorene	ND	0.23		67	65	3.0	68	62	9.2	40 - 140	30
Hexachlorobenzene	ND	0.13		64	64	0.0	62	59	5.0	40 - 140	30
Hexachlorobutadiene	ND	0.23		61	63	3.2	59	56	5.2	40 - 140	30
Hexachlorocyclopentadiene	ND	0.23		60	62	3.3	40	44	9.5	40 - 140	30
Hexachloroethane	ND	0.13		56	56	0.0	53	51	3.8	40 - 140	30
Indeno(1,2,3-cd)pyrene	ND	0.23		77	70	9.5	71	66	7.3	40 - 140	30
Isophorone	ND	0.13		54	55	1.8	57	51	11.1	40 - 140	30
Naphthalene	ND	0.23		57	58	1.7	59	55	7.0	40 - 140	30
Nitrobenzene	ND	0.13		53	58	9.0	60	53	12.4	40 - 140	30
N-Nitrosodimethylamine	ND	0.23		55	58	5.3	61	51	17.9	40 - 140	30
N-Nitrosodi-n-propylamine	ND	0.13		53	58	9.0	61	53	14.0	40 - 140	30
N-Nitrosodiphenylamine	ND	0.13		64	67	4.6	71	63	11.9	40 - 140	30
Pentachlorophenol	ND	0.23		84	84	0.0	85	74	13.8	30 - 130	30
Phenanthrene	ND	0.13		67	66	1.5	65	60	8.0	40 - 140	30
Phenol	ND	0.23		62	67	7.8	74	62	17.6	30 - 130	30
Pyrene	ND	0.23		72	70	2.8	69	61	12.3	30 - 130	30
% 2,4,6-Tribromophenol	70	%		69	68	1.5	70	64	9.0	30 - 130	30
% 2-Fluorobiphenyl	63	%		61	61	0.0	60	57	5.1	30 - 130	30
% 2-Fluorophenol	62	%		59	65	9.7	67	58	14.4	30 - 130	30
% Nitrobenzene-d5	58	%		52	55	5.6	61	54	12.2	30 - 130	30
% Phenol-d5	63	%		61	65	6.3	72	59	19.8	30 - 130	30

QA/QC Data

SDG I.D.: GCP33746

Parameter	Blank	Blk RL	QA/QC Data				SDG I.D.: GCP33746			
			LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
% Terphenyl-d14	71	%	70	68	2.9	67	60	11.0	30 - 130	30
Comment:										
Additional 8270 criteria: 20% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 15-110%, for soils 30-130%)										
QA/QC Batch 703856 (mg/Kg), QC Sample No: CP33168 (CP33748)										
<u>Volatiles - Soil (Low Level)</u>										
1,2,4-Trimethylbenzene	ND	0.001	96	110	13.6	95	84	12.3	70 - 130	30
1,3,5-Trimethylbenzene	ND	0.001	97	110	12.6	97	86	12.0	70 - 130	30
Benzene	ND	0.001	92	106	14.1	103	96	7.0	70 - 130	30
Ethylbenzene	ND	0.001	95	108	12.8	97	89	8.6	70 - 130	30
Isopropylbenzene	ND	0.001	96	109	12.7	107	96	10.8	70 - 130	30
m&p-Xylene	ND	0.002	96	109	12.7	96	88	8.7	70 - 130	30
Methyl t-butyl ether (MTBE)	ND	0.001	90	105	15.4	102	96	6.1	70 - 130	30
Naphthalene	ND	0.005	103	123	17.7	59	47	22.6	70 - 130	30
n-Butylbenzene	ND	0.001	97	111	13.5	67	59	12.7	70 - 130	30
n-Propylbenzene	ND	0.001	96	110	13.6	98	88	10.8	70 - 130	30
o-Xylene	ND	0.002	97	110	12.6	97	89	8.6	70 - 130	30
p-Isopropyltoluene	ND	0.001	97	111	13.5	80	69	14.8	70 - 130	30
sec-Butylbenzene	ND	0.001	96	109	12.7	83	73	12.8	70 - 130	30
tert-Butylbenzene	ND	0.001	95	108	12.8	95	82	14.7	70 - 130	30
Toluene	ND	0.001	91	106	15.2	99	91	8.4	70 - 130	30
% 1,2-dichlorobenzene-d4	100	%	100	100	0.0	97	97	0.0	70 - 130	30
% Bromofluorobenzene	98	%	101	101	0.0	93	94	1.1	70 - 130	30
% Dibromofluoromethane	98	%	101	101	0.0	102	102	0.0	70 - 130	30
% Toluene-d8	98	%	99	99	0.0	99	99	0.0	70 - 130	30
Comment:										
Additional 8260 criteria: 10% of LCS/LCSD compounds can be outside of acceptance criteria as long as recovery is 40-160%, 25-160% for Chloroethane-HL and Trichlorofluoromethane-HL.										
QA/QC Batch 703677 (mg/Kg), QC Sample No: CP33565 (CP33746, CP33747)										
<u>Volatiles - Soil (Low Level)</u>										
1,1,1-Trichloroethane	ND	0.005	117	120	2.5	110	112	1.8	70 - 130	30
1,1,2,2-Tetrachloroethane	ND	0.003	94	98	4.2	91	96	5.3	70 - 130	30
1,1,2-Trichloroethane	ND	0.005	105	106	0.9	88	90	2.2	70 - 130	30
1,1-Dichloroethane	ND	0.005	96	96	0.0	89	90	1.1	70 - 130	30
1,1-Dichloroethene	ND	0.005	112	110	1.8	99	99	0.0	70 - 130	30
1,2,3-Trichlorobenzene	ND	0.005	103	105	1.9	51	48	6.1	70 - 130	30
1,2,4-Trichlorobenzene	ND	0.005	99	100	1.0	56	54	3.6	70 - 130	30
1,2-Dibromo-3-chloropropane	ND	0.005	124	125	0.8	90	99	9.5	70 - 130	30
1,2-Dibromoethane	ND	0.005	105	108	2.8	84	89	5.8	70 - 130	30
1,2-Dichlorobenzene	ND	0.005	105	106	0.9	83	81	2.4	70 - 130	30
1,2-Dichloroethane	ND	0.005	116	117	0.9	101	101	0.0	70 - 130	30
1,2-Dichloropropane	ND	0.005	95	96	1.0	88	86	2.3	70 - 130	30
1,3-Dichlorobenzene	ND	0.005	99	100	1.0	87	85	2.3	70 - 130	30
1,4-Dichlorobenzene	ND	0.005	103	103	0.0	86	86	0.0	70 - 130	30
1,4-dioxane	ND	0.1	81	83	2.4	109	103	5.7	70 - 130	30
2-Hexanone	ND	0.025	79	82	3.7	15	10	40.0	70 - 130	30
4-Methyl-2-pentanone	ND	0.025	86	89	3.4	29	23	23.1	70 - 130	30
Acetone	ND	0.01	93	93	0.0	67	70	4.4	70 - 130	30
Benzene	ND	0.001	97	97	0.0	87	88	1.1	70 - 130	30
Bromochloromethane	ND	0.005	102	101	1.0	88	88	0.0	70 - 130	30
Bromodichloromethane	ND	0.005	121	122	0.8	106	107	0.9	70 - 130	30
Bromoform	ND	0.005	131	138	5.2	96	101	5.1	70 - 130	30

QA/QC Data

SDG I.D.: GCP33746

Parameter	Blank	Blk	RL	LCS	LCSD	LCS	MS	MSD	MS	%	%
				%	%	RPD	%	MSD %	MS RPD	Rec Limits	RPD Limits
Bromomethane	ND	0.005		123	121	1.6	85	78	8.6	70 - 130	30
Carbon Disulfide	ND	0.005		112	111	0.9	90	89	1.1	70 - 130	30
Carbon tetrachloride	ND	0.005		136	136	0.0	120	123	2.5	70 - 130	30
Chlorobenzene	ND	0.005		103	105	1.9	91	91	0.0	70 - 130	30
Chloroethane	ND	0.005		117	114	2.6	105	105	0.0	70 - 130	30
Chloroform	ND	0.005		104	103	1.0	94	96	2.1	70 - 130	30
Chloromethane	ND	0.005		94	89	5.5	68	70	2.9	70 - 130	30
cis-1,2-Dichloroethene	ND	0.005		99	97	2.0	87	87	0.0	70 - 130	30
cis-1,3-Dichloropropene	ND	0.005		113	114	0.9	87	88	1.1	70 - 130	30
Cyclohexane	ND	0.005		88	89	1.1	80	80	0.0	70 - 130	30
Dibromochloromethane	ND	0.003		132	132	0.0	105	110	4.7	70 - 130	30
Dichlorodifluoromethane	ND	0.005		117	111	5.3	83	86	3.6	70 - 130	30
Ethylbenzene	ND	0.001		102	103	1.0	90	91	1.1	70 - 130	30
Isopropylbenzene	ND	0.001		100	102	2.0	91	90	1.1	70 - 130	30
m&p-Xylene	ND	0.002		99	100	1.0	78	78	0.0	70 - 130	30
Methyl ethyl ketone	ND	0.005		79	83	4.9	43	41	4.8	70 - 130	30
Methyl t-butyl ether (MTBE)	ND	0.001		114	112	1.8	97	99	2.0	70 - 130	30
Methylacetate	ND	0.005		99	101	2.0	45	33	30.8	70 - 130	30
Methylcyclohexane	ND	0.005		97	99	2.0	84	83	1.2	70 - 130	30
Methylene chloride	ND	0.005		108	105	2.8	95	96	1.0	70 - 130	30
o-Xylene	ND	0.002		100	101	1.0	87	87	0.0	70 - 130	30
Styrene	ND	0.005		98	98	0.0	77	75	2.6	70 - 130	30
Tetrachloroethene	ND	0.005		111	109	1.8	102	102	0.0	70 - 130	30
Toluene	ND	0.001		104	104	0.0	93	92	1.1	70 - 130	30
trans-1,2-Dichloroethene	ND	0.005		110	108	1.8	97	98	1.0	70 - 130	30
trans-1,3-Dichloropropene	ND	0.005		126	126	0.0	92	94	2.2	70 - 130	30
Trichloroethene	ND	0.005		107	107	0.0	97	97	0.0	70 - 130	30
Trichlorofluoromethane	ND	0.005		132	130	1.5	117	121	3.4	70 - 130	30
Trichlorotrifluoroethane	ND	0.005		120	121	0.8	112	111	0.9	70 - 130	30
Vinyl chloride	ND	0.005		106	102	3.8	85	87	2.3	70 - 130	30
% 1,2-dichlorobenzene-d4	96	%		101	102	1.0	98	98	0.0	70 - 130	30
% Bromofluorobenzene	101	%		98	99	1.0	92	92	0.0	70 - 130	30
% Dibromofluoromethane	99	%		97	99	2.0	94	98	4.2	70 - 130	30
% Toluene-d8	91	%		101	100	1.0	101	100	1.0	70 - 130	30

Comment:

Additional 8260 criteria: 10% of LCS/LCSD compounds can be outside of acceptance criteria as long as recovery is 40-160%, 25-160% for Chloroethane-HL and Trichlorofluoromethane-HL.

I = This parameter is outside laboratory LCS/LCSD specified recovery limits.

m = This parameter is outside laboratory MS/MSD specified recovery limits.

r = This parameter is outside laboratory RPD specified recovery limits.

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

RPD - Relative Percent Difference

LCS - Laboratory Control Sample

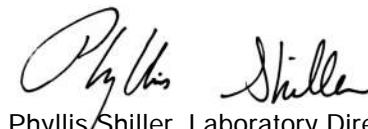
LCSD - Laboratory Control Sample Duplicate

MS - Matrix Spike

MS Dup - Matrix Spike Duplicate

NC - No Criteria

Intf - Interference



Phyllis Shiller, Laboratory Director  
November 09, 2023

Thursday, November 09, 2023

Criteria: NY: 375, 375COM, 375RRS, 375RS, CP51S

State: NY

## Sample Criteria Exceedances Report

GCP33746 - LABELLA-LATHAM

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	RL Criteria	Analysis Units
CP33747	NI-SM	Nickel	NY / 375-6.8 Metals / Unrestricted Use Soil	55.9	0.39	30	30	mg/Kg
CP33747	ZN-SM	Zinc	NY / 375-6.8 Metals / Unrestricted Use Soil	174	7.8	109	109	mg/Kg

Phoenix Laboratories does not assume responsibility for the data contained in this exceedance report. It is provided as an additional tool to identify requested criteria exceedences. All efforts are made to ensure the accuracy of the data (obtained from appropriate agencies). A lack of exceedence information does not necessarily suggest conformance to the criteria. It is ultimately the site professional's responsibility to determine appropriate compliance.



**Environmental Laboratories, Inc.**  
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045  
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## Analysis Comments

November 09, 2023

SDG I.D.: GCP33746

The following analysis comments are made regarding exceptions to criteria not already noted in the Analysis Report or QA/QC Report:

### SVOA Narration

**CHEM29 10/30/23-1:** CP33746, CP33747

For 8270 full list, the DDT breakdown and pentachlorophenol & benzidine peak tailing were evaluated in the DFTPP tune and were found to be in control.

For 8270 BN list, benzidine peak tailing was evaluated in the DFTPP tune and was found to be in control.

The following Initial Calibration compounds did not meet RSD% criteria: Bis(2-ethylhexyl)phthalate 21% (20%)

The following Initial Calibration compounds did not meet maximum RSD% criteria: None.

The following Initial Calibration compounds did not meet recommended response factors: % 2,4,6-Tribromophenol 0.043 (0.05), 2-Nitrophenol 0.041 (0.1), Bis(2-chloroethyl)ether 0.676 (0.7), Hexachlorobenzene 0.076 (0.1)

The following Initial Calibration compounds did not meet minimum response factors: % 2,4,6-Tribromophenol 0.043 (0.05), 2-Nitrophenol 0.041 (0.05)

The following Continuing Calibration compounds did not meet recommended response factors: % 2,4,6-Tribromophenol 0.046 (0.05), 2-Nitrophenol 0.045 (0.1), Bis(2-chloroethyl)ether 0.615 (0.7), Hexachlorobenzene 0.073 (0.1)

The following Continuing Calibration compounds did not meet minimum response factors: None.

Up to eight compounds can be outside of ICAL %RSD criteria and up to sixteen compounds can be outside of CCAL %Dev criteria if less than 40%.

### VOA Narration

**CHEM03 10/26/23-1:** CP33748

The following Initial Calibration compounds did not meet recommended response factors: Ethylbenzene 0.375 (0.4)

The following Initial Calibration compounds did not meet minimum response factors: None.

The following Continuing Calibration compounds did not meet recommended response factors: Ethylbenzene 0.322 (0.4)

The following Continuing Calibration compounds did not meet minimum response factors: None.

Up to eight compounds can be outside of ICAL %RSD criteria and up to sixteen compounds can be outside of CCAL %Dev criteria if less than 40%.

**CHEM18 10/25/23-2:** CP33746, CP33747

The following Initial Calibration compounds did not meet RSD% criteria: 1,2-Dibromo-3-chloropropane 26% (20%)

The following Initial Calibration compounds did not meet maximum RSD% criteria: None.

The following Initial Calibration compounds did not meet recommended response factors: Bromoform 0.091 (0.1), Dibromochloromethane 0.192 (0.2)

The following Initial Calibration compounds did not meet minimum response factors: None.

Up to eight compounds can be outside of ICAL %RSD criteria and up to sixteen compounds can be outside of CCAL %Dev criteria if less than 40%.



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## NY Temperature Narration

November 09, 2023

SDG I.D.: GCP33746

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The samples in this delivery group were received at 1.7°C.  
(Note acceptance criteria for relevant matrices is above freezing up to 6°C)

