



VIA ELECTRONIC MAIL

January 14, 2019

Mordechai Hirsch
160 Heyward Street
Brooklyn, NY 11206

Re: *Limited Phase II Investigation*
188 East 135th Street, Bronx - Tax Map No. 2-2323-13

Dear Mr. Hirsch:

Environmental Business Consultants (EBC) is pleased to provide this letter report documenting the results of the Limited Phase II Environmental Site Assessment (ESA) performed at the above-referenced property on December 12, 2018, in accordance with EBC's proposal, dated December 5, 2018.

Background

The Site consists of a single, irregular-shaped 0.57-acre parcel located on the northwestern corner of intersection between East 135th Street, which runs north to south along the eastern site perimeter, and Park Avenue, in the Mott Haven section of the Borough of the Bronx, New York City, Bronx County, New York (Figure 1). The property is currently developed with 1-story commercial building. The building occupies the east-central portion of the parcel, fronts to the north, and has a reported area of 6,500 SF. Asphalt-paved parking lots and unpaved yard areas comprise the remainder of the site. Sidewalks are located to the south and east, along Park Avenue and East 135th Street, respectively. The building is currently vacant, but was formerly occupied by the Padded Wagon (a moving company) and several other commercial tenants.

A prior Phase I Screening Report, prepared by EBC, dated in January 2018, identified several recognized environmental conditions (RECs) and other environmental concerns related to the historic use of the property, including a railroad yard, a coal yard, a contractor's storage/supply yard, its potential use as a dry cleaner, and the historic industrial use of the surrounding properties.

In addition, the property was assigned an E-designation (E-227) for Hazmat and Noise during the Lower Concourse Rezoning and Related Actions completed by the City in June 2009 (CEQR 08DCP071X). The E-designations will have to be addressed before the property can be redeveloped. The scope of this Limited Phase II Investigation represents a preliminary assessment of the property to evaluate subsurface conditions and determine if the site is eligible for inclusion in the New York State Brownfields Program, and does not fully satisfy the requirements of the E-Designation Program.



Soil Boring Investigation

To evaluate potential impacts related to the historic use of the site, EBC conducted a soil boring investigation consisting of five soil borings (SB1 through SB5) at representative locations across the site to determine if additional investigation and/or remediation is warranted. Soil boring locations are shown on Figure 2.

At each boring location soil samples were collected continuously from grade to the water table (approximately 10 to 12 feet) using a Geoprobe™. The Geoprobe™ uses direct push technology to drive core samplers to the desired depth for soil sample collection. This method can be performed quickly, so if refusal occurs, a new location can be accessed with minimal effort. Soil samples were characterized by an EBC environmental scientist and inspected for visual and olfactory evidence of contamination (i.e. staining and/or odors). Non-disposable sampling equipment was cleaned using a potable water and Alconox detergent wash followed by a potable water rinse prior to the collection of each sample. Upon collection, the samples were placed in pre-cleaned laboratory supplied glassware and stored in a cooler packed with ice for transport to the laboratory.

Soils at the site consisted generally of brown to black silty sand, with fill material (concrete, brick and asphalt) extending to a depth of 10 feet below grade. At borings SB3 and SB5, some clay and silty clay were present at depths below 8 to 10 feet. Additionally, the fill materials were very loose in some areas and sample recovery was poor. No evidence of petroleum impacts (i.e., staining or odors) was observed and no elevated PID readings were noted. Soil boring logs are included as Attachment A.

As part of the field activities, five soil samples (one per boring), either the sample exhibiting the highest degree of impact (visual or olfactory indicators) or the deepest interval above the water table, were submitted to Phoenix Environmental Laboratories, Inc., of Manchester, CT, a New York State-certified laboratory (No. 11301) for analysis. The samples submitted were from the following depths SB1 (10-12'), SB2 and SB3 (7-9') and SB4 and SB5 (8-10'). Soil samples were analyzed volatile organic compounds (VOCs) using United States Environmental Protection Agency (USEPA) Method 8260. Semi-volatile organic compounds (SVOCs) using USEPA Method 8270, and Target Analyte List (TAL) metals using USEPA Methods 6010 and 7471. These methods are specified by the New York State Department of Environmental Conservation (NYSDEC) in the evaluation of petroleum (gasoline, diesel and heating oil) spill, typical industrial solvents, and also consistent with NYCOER's E-Designation requirements.

Soil Analytical Results

Soil analytical results were compared to the NYSDEC's Division of Environmental Remediation 6 NYCRR Part 375 Soil Cleanup Objective tables (Table 375-6.8[a]: Unrestricted Use Soil Cleanup Objectives [UUSCOs]), CP-51 Soil Cleanup Guidance (Tables 2 and 3), and the Restricted Residential Use Soil Cleanup Objectives (RRUSCOs) (Part 375 Table 375-6.8[b]) to determine if additional investigation and/or remediation is warranted.

The analytical results revealed that the VOC acetone was detected in each of the five samples, at concentrations between 9.8 and 36 micrograms per kilogram (ug/kg), which are below its UUSCO of 50 ug/kg. It should be noted that acetone is a common laboratory contaminant, and its presence in these samples may be attributable to laboratory cross-contamination. Naphthalene was also detected in sample SB3, at a concentration of 71 ug/kg. There is no UUSCO for this compound. No other VOCs were detected in any of the samples at concentrations exceeding their respective laboratory method detection limits (MDLs).

The SVOCs benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, dibenzo(a,h)anthracene, and indeno(1,2,3-cd)pyrene, were detected in soil sample SB1 (10-12') at concentrations exceeding their respective UUSCOs. In addition, the concentrations of these compounds, except for benzo(k)fluoranthene and chrysene, also exceeded their respective RRUSCOs. Twelve additional SVOCs were also detected in sample SB1 (10-12'), but at concentrations below their UUSCOs. One or more SVOCs were also detected in the samples from borings SB2, SB3 and SB4, but at concentrations below their UUSCOs. No SVOCs were detected in the soil sample from boring SB5 at concentrations above their respective laboratory MDLs.

The metals arsenic, cadmium, chromium, copper, lead, mercury, nickel and zinc, were detected in one of more of the samples SB1 through SB4 at concentrations exceeding their respective UUSCOs. The concentration of cadmium (SB1 and SB2), copper (SB1), lead (SB1, SB2, and SB3), and mercury (SB1 through SB4) also exceeded their respective RRUSCOs. Several additional metals were also detected in samples SB1 through SB4 and in sample SB5, but at concentrations below their respective UUSCOs.

Laboratory analytical results for the soil samples are summarized on Tables 1 through 3, and the laboratory reports are provided as Attachment B.

Groundwater Sampling

As groundwater was encountered at approximately 10 to 12 feet below grade, three of the five borings (SB1, SB3 and SB4) were extended to a depth three feet below the water table to facilitate the collection of groundwater samples (GW1 through GW3). Groundwater samples were collected using the Geoprobe™ equipped with a four-foot long mill slot sampler or equivalent. The sampler was first be driven to the desired depth (approximately three feet below the water table). This allows the sampler screen to intersect the water table and allow floating product or petroleum sheens (if present) to be documented. A piece of disposable polyethylene tubing with a stainless steel check valve was then inserted through the probe rods into the water bearing zone and the tubing hand oscillated to obtain the sample. The groundwater samples were collected directly from the tubing into pre-cleaned laboratory supplied glassware and stored in a cooler packed with ice for transport to the laboratory. Non-disposable sampling equipment was cleaned using a distilled water and Alconox detergent wash followed by a potable water rinse prior to the collection of each sample. The groundwater samples were analyzed for VOCs by EPA Method 8260. Groundwater sample locations are shown on Figure 2.

Groundwater Analytical Results

Groundwater analytical results were compared to New York State 6NYCRR Part 703.5 Class GA groundwater quality standards (GQS).

The analytical results revealed that 10 VOCs, primarily petroleum-related compounds detected in sample GW1, which was collected from boring SB1, but at concentrations below their respective groundwater standards. Two VOCs were also detected in sample GW3, collected from boring SB4, but at concentrations below their respective groundwater standards. No VOCs were detected in sample GW2 at concentrations exceeding their respective laboratory method detection limits (MDLs).

Laboratory analytical results for the groundwater samples are summarized on Table 4, and the laboratory reports are provided as Attachment B.

Soil Vapor Sampling

To evaluate potential migration of VOCs associated with historic industrial use of the site and their potential impact to soil vapor beneath the building at the site. EBC conducted a soil vapor survey, which consisted of the collection and analysis of two sub-surface soil vapor samples (SV1 and SV2) at

representative locations and one sub-slab soil vapor sample (SS1) from beneath the central portion of the site building. At both of the subsurface soil vapor sample locations, temporary vapor points consisting of six-inch long stainless steel screen point samplers and polyethylene tubing, were installed to a depth of approximately 6 feet below grade (midpoint between the ground surface and the water table) using Geoprobe™ equipment. The sub-slab vapor point also consisted of a six-inch long stainless steel screen point sampler and polyethylene tubing, installed to a depth of approximately one foot below the floor slab using a Bosch rotary hammer drill to penetrate the floor and drill to the desired depth.

At each location, the annular space between the polyethylene tubing and the boreholes was then backfilled to grade with sand, with a bentonite clay seal emplaced above the sand at the ground surface to prevent ambient air from being drawn into the borehole and mixing with the soil vapor to be sampled. The above-grade end of the tubing was then attached to a hand pump and ambient air within the tubing was purged to ensure the collection of a representative sample. The tubing was then attached directly to a 6-liter laboratory-supplied SUMMA vacuum canister, equipped with laboratory calibrated flow controllers. The air (soil vapor) samples were collected for a period of approximately two hours at a rate of 0.05 liter per minute to obtain the required sample volume. After collection, the canisters were properly labeled and shipped under chain-of-custody to Phoenix Environmental Laboratories for analysis of VOCs using USEPA Method TO-15.

Soil Vapor Analytical Results

Soil vapor analytical results were compared to the compounds listed in Table 3.1 of the Air Guideline Values Derived by the NYSDOH located in the New York State Department of Health (NYSDOH) Final Guidance for Evaluating Soil Vapor Intrusion dated October 2006.

Each of the three soil vapor/sub-slab samples contained tetrachloroethene (PCE) at concentrations (0.62 µg/m³ to 5.73 µg/m³), within the mitigation range established within the State DOH soil vapor guidance matrix. The chlorinated VOCs trichloroethene (max. 0.35 µg/m³), vinyl chloride (max. 0.49 µg/m³) and carbon tetrachloride (max. 0.33 µg/m³) were also detected in one or more of the three soil vapor samples, but at concentrations below applicable criteria. The chlorinated VOCs 1,1-dichloroethene and cis-1,2-dichloroethene were not detected in the soil vapor/sub-slab samples. Low levels of petroleum-related VOCs were also present in each of the samples. The total concentration of petroleum-related VOCs (Total BTEX) ranged from 29.45 µg/m³ in SV2 to 57.71 µg/m³ in SS1. Individual BTEX compounds included benzene (5.87 µg/m³), ethylbenzene (5.47 µg/m³), toluene (max. 21.6 µg/m³), m&p xylenes (max. 18.3 µg/m³) and o-xylene (max. 6.47 µg/m³). Soil vapor/sub-slab sample locations are shown on Figure 3.

Laboratory analytical results for the soil vapor samples are summarized on Table 5, and the laboratory reports are provided as Attachment B.

Conclusions and Recommendations

Soil samples were collected from five representative soil borings drilled across the site and analyzed for the presence of VOCs, SVOCs and metals. Analytical results indicate the presence of several SVOCs at concentrations exceeding applicable regulatory criteria in one sample (SB1 10-12') at the northeastern portion of the site. No evidence of petroleum staining or odors was observed in this boring. One or more heavy metals were detected in four of the five soil samples at concentrations exceed either their UUSCO and/or RRUSCO. The presence of these compounds (SVOCs and metals) is consistent with typical urban fill materials and not indicative of a petroleum release.

Several VOCs were detected in two of the three groundwater samples, but at concentrations below applicable regulatory criteria.



Soil vapor and sub-slab soil vapor samples contained low concentrations of chlorinated and petroleum-related compounds, none of which exceeded applicable regulatory criteria. As such, there is no evidence of a soil vapor intrusion condition at the site.

In general, the findings of the Phase II investigation are consistent with the presumed conditions outlined in our November 20, 2018 environmental review letter, which stated that the site was likely underlain by typical urban fill materials impacted by metals and other contaminants. In this case SVOCs in the northeastern portions. Because the property has been assigned an E-designation for hazardous materials by New York City, a more thorough subsurface will be required prior to any redevelopment of the site, and any soils removed for construction will have to be properly disposed of. In addition, to the soil/fill material disposal, NYCOER will likely require that any new building be constructed with a sub-slab vapor barrier, although, based upon the soil vapor sample results and the absence of VOCs in site soils and groundwater the installation of a sub-slab depressurization system is unlikely to be warranted.

Finally, as noted in our prior environmental review letter, the property is located within an Enzone, as well as designated Brownfield Opportunity Area. Based upon the levels of SVOCs and metals detected in onsite fill materials, you could opt into the State Brownfield program, which would provide tax credits for the clean-up and the new building.

We appreciate the opportunity to assist you with this project. Should you have any questions or comments, please do not hesitate to contact me.

Very truly yours,

Environmental Business Consultants

Keith W. Butler
Senior Project Manager



TABLES

TABLE 1
188 East 135th Street
Bronx, New York
Soil Analytical Results
Volatile Organic Compounds

COMPOUND	NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives*	NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives*	SB1		SB2		SB3		SB4		SB5	
			(10-12')		(7-9')		(7-9')		(8-10')		(8-10')	
			12/12/2018 µg/Kg	RL	12/12/2018 µg/Kg	RL	12/12/2018 µg/Kg	RL	12/12/2018 µg/Kg	RL	12/12/2018 µg/Kg	RL
1,1,1,2-Tetrachloroethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,1,1-Trichloroethane	680	100,000	< 19	19	< 4.8	4.8	< 25	25	< 5.7	5.7	< 5.0	5.0
1,1,2-Tetrachloroethane			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,1,2-Trichloroethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,1-Dichloroethane	270	26,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,1-Dichloroethane	330	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,1-Dichloropropene			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,2,3-Trichlorobenzene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2,3-Trichloropropane			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2,4-Trichlorobenzene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2,4-Trimethylbenzene	3,600	52,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2-Dibromo-3-chloropropane			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2-Dibromomethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,2-Dichlorobenzene	1,100	100,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,2-Dichloroethane	20	3,100	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,2-Dichloroethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,3,5-Trimethylbenzene	8,400	52,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,3-Dichlorobenzene	2,400	4,900	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,3-Dichloropropane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
1,4-Dichlorobenzene	1,800	13,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
1,4-dioxane	100	13,000	< 72	72	< 72	72	< 94	94	< 95	95	< 75	75
2,2-Dichloropropane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
2-Chlorotoluene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
2-Hexanone (Methyl Butyl Ketone)			< 24	24	< 24	24	< 31	31	< 28	28	< 25	25
2-Isopropyltoluene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
4-Chlorotoluene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
4-Methyl-2-Pentanone			< 24	24	< 24	24	< 31	31	< 28	28	< 25	25
Acetone	50	100,000	9.8	24	23	24	23	31	10	28	36	25
Acrolein			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Acrylonitrile			< 19	19	< 9.6	9.6	< 13	13	< 11	11	< 20	20
Benzene	60	4,800	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Bromobenzene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
Bromochloromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Bromodichloromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Bromoform			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Bromomethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Carbon Disulfide			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Carbon tetrachloride	760	2,400	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Chlorobenzene	1,100	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Chloroethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Chloroform	370	49,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Chloromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
cis-1,2-Dichloroethene	250	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
cis-1,3-Dichloropropane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Dibromochloromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Dibromomethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Dichlorodifluoromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Ethylbenzene	1,000	41,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Hexachlorobutadiene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
Isopropylbenzene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
m&p-Xylenes	260	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Methyl Ethyl Ketone (2-Butanone)	120	100,000	< 29	29	< 29	29	< 38	38	< 34	34	< 30	30
Methyl t-butyl ether (MTBE)	930	100,000	< 9.7	9.7	< 9.6	9.6	< 13	13	< 11	11	< 10	10
Methylene chloride	50	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Naphthalene			< 320	320	< 4.8	4.8	71	320	< 5.7	5.7	< 5.0	5.0
n-Butylbenzene	12,000	100,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
n-Propylbenzene	3,900	100,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
o-Xylene	260	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
p-Isopropyltoluene			< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
sec-Butylbenzene	11,000	100,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
Styrene			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
tert-butyl alcohol			< 97	97	< 96	96	< 130	130	< 110	110	< 100	100
tert-Butylbenzene	5,900	100,000	< 320	320	< 4.8	4.8	< 320	320	< 5.7	5.7	< 5.0	5.0
Tetrachloroethene	1,300	19,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Tetrahydrofuran (THF)			< 9.7	9.7	< 9.6	9.6	< 13	13	< 11	11	< 10	10
Toluene	700	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
trans-1,2-Dichloroethene	190	100,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
trans-1,3-Dichloropropane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
trans-1,4-dichloro-2-butene			< 640	640	< 9.6	9.6	< 640	640	< 11	11	< 10	10
Trichloroethene	470	21,000	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Trichlorofluoromethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Trichlorotrifluoroethane			< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Vinyl Chloride	20	900	< 4.8	4.8	< 4.8	4.8	< 6.3	6.3	< 5.7	5.7	< 5.0	5.0
Total BTEX Concentration			0		0		0		0		0	
Total VOCs Concentration			9.8		23		94		10		36	

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

BCG - Below Cellar Grade

RL- Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSOC Guidance Value

TABLE 2
188 East 135th Street
Bronx, New York
Soil Analytical Results
Semi-Volatile Organic Compounds

COMPOUND	NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives*	NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives*	SB1		SB2		SB3		SB4		SB5	
			(10-12) 12/12/2018		(7-9) 12/12/2018		(7-9) 12/12/2018		(8-10) 12/12/2018		(8-10) 12/12/2018	
			Result	RL	Result	RL	Result	RL	Result	RL	Result	RL
1,2,4,5-Tetrachlorobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
1,2,4-Trichlorobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
1,2-Dichlorobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
1,2-Diphenylhydrazine			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
1,3-Dichlorobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
1,4-Dichlorobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2,4,5-Trichlorophenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2,4,6-Trichlorophenol			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
2,4-Dichlorophenol			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
2,4-Dimethylphenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2,4-Dinitrophenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2,4-Dinitrotoluene			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
2,6-Dinitrotoluene			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
2-Chloronaphthalene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2-Chlorophenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2-Methylnaphthalene			250	260	< 270	270	< 260	260	< 250	250	< 270	270
2-Methylphenol (o-cresol)	330	100,000	< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2-Nitroaniline			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
2-Nitrophenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
3,4-Methylphenol (m&p-cresol)	330	100,000	< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
3,3'-Dichlorobenzidine			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
3-Nitroaniline			< 380	380	< 390	390	< 380	380	< 360	360	< 390	390
4,6-Dinitro-2-methylphenol			< 230	230	< 230	230	< 230	230	< 220	220	< 230	230
4-Bromophenyl phenyl ether			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
4-Chloro-3-methylphenol			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
4-Chloroaniline			< 300	300	< 310	310	< 300	300	< 290	290	< 310	310
4-Chlorophenyl phenyl ether			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
4-Nitroaniline			< 380	380	< 390	390	< 380	380	< 360	360	< 390	390
4-Nitrophenol			< 380	380	< 390	390	< 380	380	< 360	360	< 390	390
Acenaphthene	20,000	100,000	490	260	< 270	270	< 260	260	< 250	250	< 270	270
Acenaphthylene	100,000	100,000	550	260	< 270	270	< 260	260	< 250	250	< 270	270
Acetophenone			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Aniline			< 300	300	< 310	310	< 300	300	< 290	290	< 310	310
Anthracene	100,000	100,000	1,500	260	< 270	270	140	260	380	250	< 270	270
Benz(a)anthracene	1,000	1,000	3,400	260	< 270	270	430	260	700	250	< 270	270
Benzidine			< 380	380	< 390	390	< 380	380	< 360	360	< 390	390
Benzof(a)pyrene	1,000	1,000	3,200	190	< 190	190	470	190	640	180	< 200	200
Benzo(b)fluoranthene	1,000	1,000	3,000	260	< 270	270	400	260	530	250	< 270	270
Benzo(ghi)perylene	100,000	100,000	1,800	260	< 270	270	330	260	330	250	< 270	270
Benzo(k)fluoranthene	800	3,900	2,700	260	< 270	270	380	260	500	250	< 270	270
Benzoic acid			< 1900	1,900	< 1900	1,900	< 1900	1,900	< 1800	1,800	< 2000	2,000
Benzyl butyl phthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Bis(2-chloroethoxy)methane			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Bis(2-chloroethoxy)ether			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
Bis(2-chloroisopropyl)ether			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Bis(2-ethylhexyl)phthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Carbazole			360	190	< 190	190	< 190	190	< 180	180	< 200	200
Chrysene	1,000	3,900	3,400	260	< 270	270	450	260	680	250	< 270	270
Dibenz(a,h)anthracene	330	330	720	190	< 190	190	< 190	190	< 180	180	< 200	200
Dibenzofuran	7,000	59,000	320	260	< 270	270	< 260	260	< 250	250	< 270	270
Diethyl phthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Dimethylphthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Di-n-butylphthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Di-n-octylphthalate			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Fluoranthene	100,000	100,000	7,000	260	320	270	780	260	1,700	250	< 270	270
Fluorene	30,000	100,000	470	260	< 270	270	< 260	260	< 250	250	< 270	270
Hexachlorobenzene			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
Hexachlorobutadiene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Hexachlorocyclopentadiene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Hexachloroethane			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
Indeno(1,2,3-cd)pyrene	500	500	2,100	260	< 270	270	330	260	360	250	< 270	270
Isophorone			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
Naphthalene	12,000	100,000	210	260	< 270	270	< 260	260	< 250	250	< 270	270
Nitrobenzene			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
N-Nitrosodimethylamine			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
N-Nitrosodi-n-propylamine			< 190	190	< 190	190	< 190	190	< 180	180	< 200	200
N-Nitrosodiphenylamine			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Pentachloronitrobenzene			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Pentachlorophenol	800	6,700	< 230	230	< 230	230	< 230	230	< 220	220	< 230	230
Phenanthrene	100,000	100,000	5,000	260	300	270	520	260	1,300	250	< 270	270
Phenol	330	100,000	< 260	260	< 270	270	< 260	260	< 250	250	< 270	270
Pyrene	100,000	100,000	6,500	260	< 270	270	740	260	1,500	250	< 270	270
Pyridine			< 260	260	< 270	270	< 260	260	< 250	250	< 270	270

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

BCG - Below Cellar Grade

RL - Reporting Limit

Bold/highlighted - Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted - Indicated exceedance of the NYSDEC RRSO Guidance Value

TABLE 3
188 East 135th Street
Bronx, New York
Soil Analytical Results
Metals

COMPOUND	NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives*	NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives*	SB1		SB2		SB3		SB4		SB5	
			(10-12') 12/12/2018 µg/Kg		(7-9') 12/12/2018 µg/Kg		(7-9') 12/12/2018 µg/Kg		(8-10') 12/12/2018 µg/Kg		(8-10') 12/12/2018 µg/Kg	
			Result	RL	Result	RL	Result	RL	Result	RL	Result	RL
Aluminum			5,850	37	9,320	40	9,190	37	9,910	38	12,400	37
Antimony			15.1	3.7	< 4.0	4.0	< 3.7	3.7	< 3.8	3.8	< 3.7	3.7
Arsenic	13	16	15.8	0.75	8.39	0.80	8.23	0.74	5.21	0.76	2.64	0.74
Barium	350	350	168	0.7	285	0.8	203	0.7	71.1	0.8	27.3	0.7
Beryllium	7.2	14	0.44	0.30	0.52	0.32	0.56	0.29	0.49	0.30	0.53	0.29
Cadmium	2.5	2.5	3.44	0.37	2.61	0.40	0.9	0.37	0.43	0.38	< 0.37	0.37
Calcium			48,200	37	44,500	40	6,280	3.7	1,540	3.8	7,670	3.7
Chromium	30	180	46.3	0.37	23.2	0.40	34.8	0.37	17.4	0.38	17.9	0.37
Cobalt			10.3	0.37	9.08	0.40	8.51	0.37	10.7	0.38	6.1	0.37
Copper	50	270	308	7.5	159	0.8	63.9	0.7	26.8	0.8	25.3	0.7
Iron			59,600	37	39,300	40	25,000	37	20,000	38	14,700	37
Lead	63	400	662	7.5	959	8.0	382	7.4	62.6	0.8	11.7	0.7
Magnesium			14,000	37	4,420	4.0	3,000	3.7	2,980	3.8	6,300	37
Manganese	1,600	2,000	597	3.7	398	4.0	228	3.7	293	3.8	110	0.37
Mercury	0.18	0.81	1.29	0.15	13	1.4	0.53	0.14	0.24	0.13	0.03	0.03
Nickel	30	140	38.3	0.37	18.5	0.40	43.2	0.37	16.3	0.38	14.4	0.37
Potassium			1,120	7	1,840	8	1,310	7	2,030	8	1,470	7
Selenium	3.9	36	< 1.5	1.5	< 1.6	1.6	< 1.5	1.5	< 1.5	1.5	< 1.5	1.5
Silver	2	36	< 0.37	0.37	< 0.40	0.40	< 0.37	0.37	< 0.38	0.38	< 0.37	0.37
Sodium			494	7	396	8	375	7	125	8	289	7
Thallium			< 1.5	1.5	< 1.6	1.6	< 1.5	1.5	< 1.5	1.5	< 1.5	1.5
Vanadium			56.3	0.37	25.3	0.40	23.5	0.37	26.3	0.38	22.5	0.37
Zinc	109	2,200	720	7.5	877	8.0	195	7.4	85.6	0.8	64.4	0.7

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

BCG - Below Cellar Grade

RL- Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

TABLE 4
188 East 135th Street
Bronx, New York
Ground Water Analytical Results
Volatile Organic Compounds

Compound	NYSDEC Groundwater Quality Standards µg/L	GW1		GW2		GW3	
		12/12/2018		12/12/2018		12/12/2018	
		Results	RL	Results	RL	Results	RL
1,1,1,2-Tetrachloroethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,1,1-Trichloroethane	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
1,1,2,2-Tetrachloroethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,1,2-Trichloroethane	1	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,1-Dichloroethane	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
1,1-Dichloroethene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,1-Dichloropropene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,2,3-Trichlorobenzene		< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,2,3-Trichloropropane	0.04	< 0.25	0.25	< 0.25	0.25	< 0.25	0.25
1,2,4-Trichlorobenzene		< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,2,4-Trimethylbenzene	5	1.8	1.0	< 1.0	1.0	< 1.0	1.0
1,2-Dibromo-3-chloropropane	0.04	< 0.50	0.50	< 0.50	0.50	< 0.50	0.50
1,2-Dibromoethane	0.0006	< 0.25	0.25	< 0.25	0.25	< 0.25	0.25
1,2-Dichlorobenzene		< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,2-Dichloroethane	0.6	< 0.60	0.60	< 0.60	0.60	< 0.60	0.60
1,2-Dichloropropane	1	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,3,5-Trimethylbenzene	5	0.34	1.0	< 1.0	1.0	< 1.0	1.0
1,3-Dichlorobenzene	3	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,3-Dichloropropane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,4-Dichlorobenzene		< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
1,4-dioxane		< 100	100	< 100	100	< 100	100
2,2-Dichloropropane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
2-Chlorotoluene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
2-Hexanone	50	< 2.5	2.5	< 2.5	2.5	< 2.5	2.5
2-Isopropyltoluene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
4-Chlorotoluene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
4-Methyl-2-pentanone		< 2.5	2.5	< 2.5	2.5	< 2.5	2.5
Acetone	50	3.9	5.0	< 5.0	5.0	3.7	5.0
Acrolein	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Acrylonitrile	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Benzene	1	< 0.70	0.70	< 0.70	0.70	< 0.70	0.70
Bromobenzene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Bromochloromethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Bromodichloromethane	50	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Bromoform	50	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Bromomethane	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Carbon Disulfide		0.4	1.0	< 1.0	1.0	0.39	1.0
Carbon tetrachloride	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Chlorobenzene	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Chloroethane	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Chloroform	7	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Chloromethane	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
cis-1,2-Dichloroethene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
cis-1,3-Dichloropropene	0.4	< 0.40	0.40	< 0.40	0.40	< 0.40	0.40
Dibromochloromethane	50	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Dibromomethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Dichlorodifluoromethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Ethylbenzene	5	0.32	1.0	< 1.0	1.0	< 1.0	1.0
Hexachlorobutadiene	0.5	< 0.50	0.50	< 0.50	0.50	< 0.50	0.50
Isopropylbenzene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
m&p-Xylene		1.2	1.0	< 1.0	1.0	< 1.0	1.0
Methyl ethyl ketone	50	< 2.5	2.5	< 2.5	2.5	< 2.5	2.5
Methyl t-butyl ether (MTBE)		< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Methylene chloride	5	< 3.0	3.0	< 3.0	3.0	< 3.0	3.0
Naphthalene	10	1.6	1.0	< 1.0	1.0	< 1.0	1.0
n-Butylbenzene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
n-Propylbenzene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
o-Xylene	5	0.6	1.0	< 1.0	1.0	< 1.0	1.0
p-Isopropyltoluene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
sec-Butylbenzene	5	1.5	1.0	< 1.0	1.0	< 1.0	1.0
Styrene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Tert-butyl alcohol		< 50	50	< 50	50	< 50	50
tert-Butylbenzene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Tetrachloroethene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Tetrahydrofuran (THF)	50	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
Toluene	5	1.3	1.0	< 1.0	1.0	< 1.0	1.0
trans-1,2-Dichloroethene	5	< 5.0	5.0	< 5.0	5.0	< 5.0	5.0
trans-1,3-Dichloropropene	0.4	< 0.40	0.40	< 0.40	0.40	< 0.40	0.40
trans-1,4-dichloro-2-butene	5	< 2.5	2.5	< 2.5	2.5	< 2.5	2.5
Trichloroethene	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Trichlorofluoromethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Trichlorotrifluoroethane	5	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0
Vinyl chloride	2	< 1.0	1.0	< 1.0	1.0	< 1.0	1.0

Notes:

RL- Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

TABLE 5
188 East 135th Street
Bronx, New York
Soil Gas - Volatile Organic Compounds

COMPOUNDS	NYSDOH Maximum Sub-Slab Value ($\mu\text{g}/\text{m}^3$) ^(a)	NYSDOH Soil Outdoor Background Levels ($\mu\text{g}/\text{m}^3$) ^(b)	SV1		SV2		SS1	
			12/12/2018 ($\mu\text{g}/\text{m}^3$)		12/12/2018 ($\mu\text{g}/\text{m}^3$)		12/12/2018 ($\mu\text{g}/\text{m}^3$)	
			Result	RL	Result	RL	Result	RL
1,1,1,2-Tetrachloroethane			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,1,1-Trichloroethane	100	<2.0 - 2.8	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,1,2,2-Tetrachloroethane		<1.5	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,1,2-Trichloroethane		<1.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,1-Dichloroethane		<1.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,1-Dichloroethene		<1.0	< 0.20	0.20	< 0.20	0.20	< 0.20	0.20
1,2,4-Trichlorobenzene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,2,4-Trimethylbenzene		<1.0	2.9	1.00	2.52	1.00	4.81	1.00
1,2-Dibromoethane		<1.5	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,2-Dichlorobenzene		<2.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,2-Dichloroethane		<1.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,2-Dichloropropane			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,2-Dichlorotetrafluoroethane			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,3,5-Trimethylbenzene		<1.0	1.41	1.00	< 1.00	1.00	1.72	1.00
1,3-Butadiene		NA	< 1.00	1.00	8.89	1.00	< 1.00	1.00
1,3-Dichlorobenzene		<2.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,4-Dichlorobenzene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
1,4-Dioxane			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
2-Hexanone			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
4-Ethyltoluene		NA	4.31	1.00	3.75	1.00	6.14	1.00
4-Isopropyltoluene			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
4-Methyl-2-pentanone			< 1.00	1.00	2.28	1.00	1.44	1.00
Acetone		NA	287	5.01	247	5.01	148	5.01
Acrylonitrile			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Benzene		<1.6 - 4.7	5.33	1.00	2.77	1.00	5.87	1.00
Benzyl Chloride		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Bromodichloromethane		<5.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Bromoform		<1.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Bromomethane		<1.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Carbon Disulfide		NA	14.7	1.00	5.13	1.00	3.64	1.00
Carbon Tetrachloride	5	<3.1	< 0.20	0.20	< 0.20	0.20	0.33	0.20
Chlorobenzene		<2.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Chloroethane		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Chloroform		<2.4	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Chloromethane		<1.0 - 1.4	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
cis-1,2-Dichloroethene		<1.0	< 0.20	0.20	< 0.20	0.20	< 0.20	0.20
cis-1,3-Dichloropropene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Cyclohexane		NA	8.57	1.00	10.1	1.00	1.07	1.00
Dibromochloromethane		<5.0	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Dichlorodifluoromethane		NA	1.46	1.00	1.69	1.00	1.5	1.00
Ethanol			27.5	1.00	42	1.00	19.8	1.00
Ethyl Acetate		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Ethylbenzene		<4.3	4.02	1.00	2.5	1.00	5.47	1.00
Heptane		NA	41.8	1.00	3.83	1.00	3.26	1.00
Hexachlorobutadiene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Hexane		<1.5	77.5	1.00	5.74	1.00	1.32	1.00
Isopropylalcohol		NA	< 1.00	1.00	1.74	1.00	2.87	1.00
Isopropylbenzene			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Xylene (m&p)		<4.3	15.1	1.00	10.3	1.00	18.3	1.00
Methyl Ethyl Ketone			22	1.00	11.5	1.00	12.5	1.00
MTBE		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Methylene Chloride		<3.4	64.9	3.00	12.2	3.00	< 3.00	3.00
n-Butylbenzene			< 1.00	1.00	< 1.00	1.00	1.22	1.00
Xylene (o)		<4.3	5.34	1.00	3.38	1.00	6.47	1.00
Propylene		NA	< 1.00	1.00	168	5.01	2.79	1.00
sec-Butylbenzene			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Styrene		<1.0	1.53	1.00	1.19	1.00	2.88	1.00
Tetrachloroethene	30		3.75	0.25	0.62	0.25	5.73	0.25
Tetrahydrofuran		NA	< 1.00	1.00	7.93	1.00	< 1.00	1.00
Toluene		1.0 - 6.1	17.5	1.00	10.5	1.00	21.6	1.00
trans-1,2-Dichloroethene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
trans-1,3-Dichloropropene		NA	< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Trichloroethene	2	<1.7	< 0.20	0.20	< 0.20	0.20	0.35	0.20
Trichlorofluoromethane		NA	< 1.00	1.00	< 1.00	1.00	1.85	1.00
Trichlorotrifluoroethane			< 1.00	1.00	< 1.00	1.00	< 1.00	1.00
Vinyl Chloride		<1.0	< 0.20	0.20	0.49	0.20	< 0.20	0.20
BTEX			47.29		29.45		57.71	
CVOCs			3.75		1.11		6.41	
Total VOCs			606.62		566.05		280.93	

Notes:

NA No guidance value or standard available

(a) Final Guidance for Evaluating Soil Vapor Intrusion in the State of New York, October 2006, New York State Department of Health.

(b) NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York, February 2005, Summary of Background Levels for Selected Compounds (NYSDOH Database, Outdoor values)

FIGURES

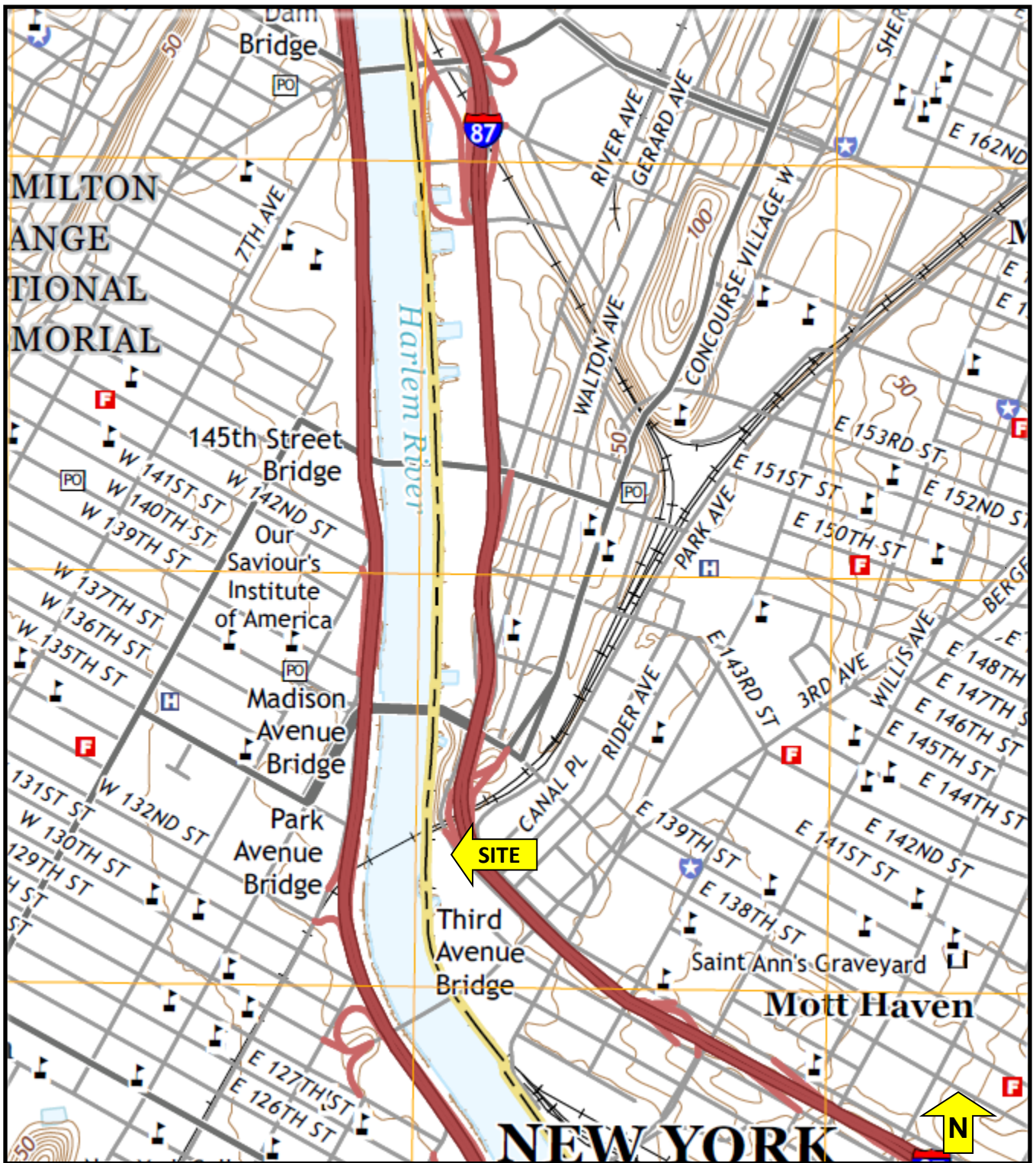


FIGURE 1 – SITE LOCATION MAP



Phone 631.504.6000
 Fax 631.924.2870

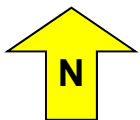
ENVIRONMENTAL BUSINESS CONSULTANTS

SITE NAME: Commercial Property
STREET ADDRESS: 188 East 135th Street
MUNICIPALITY, STATE, ZIP: Bronx, NY10451

Source: USGS



FIGURE 2 - SOIL BORING AND GROUNDWATER SAMPLE LOCATION MAP



SITE NAME: Commercial Property
STREET ADDRESS: 188 East 135th Street
MUNICIPALITY, STATE, ZIP: Bronx, NY 10036

Source: Google Earth – June 2011



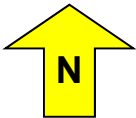
Phone 631.504.6000
 Fax 631.924.2870

ENVIRONMENTAL BUSINESS CONSULTANTS



Google Earth
©2011 B. Google

SOIL VAPOR AND SUB-SLAB SOIL VAPOR SAMPLE LOCATION MAP



SITE NAME: Commercial Property
STREET ADDRESS: 188 East 135th Street
MUNICIPALITY, STATE, ZIP: Bronx, NY 10036

Source: Google Earth – June 2011



Phone 631.504.6000
Fax 631.924.2870

ENVIRONMENTAL BUSINESS CONSULTANTS

ATTACHMENT A
SOIL BORING LOGS

ATTACHMENT B

LABORATORY ANALYTICAL REPORTS



Wednesday, December 19, 2018

Attn: Mr. Charles B. Sosik, P.G.
Environmental Business Consultants
1808 Middle Country Rd
Ridge NY 11961-2406

Project ID: 188 EAST 135TH ST BRONX NY
Sample ID#s: CC13976 - CC13983

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 have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext. 200.

Sincerely yours,

A handwritten signature in black ink that reads "Phyllis Shiller". The signature is written in a cursive style.

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
UT Lab Registration #CT00007
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

December 19, 2018

SDG I.D.: GCC13976

8260 Volatile Organics:

1,2-Dibromoethane, 1,2,3 Trichloropropane, and 1,2-Dibromo-3-chloropropane do not meet NY TOGS GA criteria, these compounds are analyzed by GC/FID method 504 or 8011 to achieve this criteria.

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

Analysis Report
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

8:15
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13976

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: SB 1 (10-12')

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.37	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Aluminum	5850	37		mg/Kg	10	12/15/18	CPP	SW6010C
Arsenic	15.8	0.75		mg/Kg	1	12/15/18	CPP	SW6010C
Barium	168	0.7		mg/Kg	1	12/15/18	EK	SW6010C
Beryllium	0.44	0.30		mg/Kg	1	12/15/18	CPP	SW6010C
Calcium	48200	37		mg/Kg	10	12/15/18	CPP	SW6010C
Cadmium	3.44	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Cobalt	10.3	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Chromium	46.3	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Copper	308	7.5		mg/kg	10	12/15/18	CPP	SW6010C
Iron	59600	37		mg/Kg	10	12/15/18	CPP	SW6010C
Mercury	1.29	0.15		mg/Kg	1	12/14/18	RS	SW7471B
Potassium	1120	7		mg/Kg	1	12/15/18	EK	SW6010C
Magnesium	14000	37		mg/Kg	10	12/15/18	CPP	SW6010C
Manganese	597	3.7		mg/Kg	10	12/15/18	CPP	SW6010C
Sodium	494	7		mg/Kg	1	12/15/18	CPP	SW6010C
Nickel	38.3	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Lead	662	7.5		mg/Kg	10	12/15/18	CPP	SW6010C
Antimony	15.1	3.7		mg/Kg	1	12/15/18	EK	SW6010C
Selenium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Thallium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Vanadium	56.3	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Zinc	720	7.5		mg/Kg	10	12/15/18	CPP	SW6010C
Percent Solid	87			%		12/13/18	AK	SW846-%Solid
Soil Extraction for SVOA	Completed					12/13/18	JJ/CK	SW3545A
Mercury Digestion	Completed					12/14/18	EV/EV	SW7471B
Total Metals Digest	Completed					12/14/18	M/AG	SW3050B

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,1-Trichloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2,2-Tetrachloroethane	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,1,2-Trichloroethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichlorobenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,3-Trichloropropane	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,4-Trichlorobenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,4-Trimethylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dibromo-3-chloropropane	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dibromoethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dichloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloropropane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
1,3,5-Trimethylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,3-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,3-Dichloropropane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
1,4-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
2,2-Dichloropropane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
2-Chlorotoluene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
2-Hexanone	ND	24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
2-Isopropyltoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
4-Chlorotoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
4-Methyl-2-pentanone	ND	24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Acetone	9.8	JS 24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	9.7	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Benzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Bromobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Bromochloromethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Bromodichloromethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Bromoform	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Bromomethane	ND	4.8	1.9	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon Disulfide	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon tetrachloride	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Chlorobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroform	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloromethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,2-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,3-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromochloromethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromomethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Dichlorodifluoromethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Ethylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Hexachlorobutadiene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Isopropylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
m&p-Xylene	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl Ethyl Ketone	ND	29	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl t-butyl ether (MTBE)	ND	9.7	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Methylene chloride	ND	4.8	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Naphthalene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
n-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
n-Propylbenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
o-Xylene	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
p-Isopropyltoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
sec-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Styrene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
tert-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Tetrachloroethene	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrahydrofuran (THF)	ND	9.7	2.4	ug/Kg	1	12/15/18	JLI	SW8260C
Toluene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,2-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,3-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,4-dichloro-2-butene	ND	640	160	ug/Kg	50	12/15/18	JLI	SW8260C
Trichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorofluoromethane	ND	4.8	0.97	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorotrifluoroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Vinyl chloride	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	100			%	50	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	99			%	50	12/15/18	JLI	70 - 130 %
% Dibromofluoromethane	100			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	94			%	1	12/15/18	JLI	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	72		ug/kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	121			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	73			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	94			%	1	12/15/18	JLI	70 - 130 %
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	19		ug/Kg	1	12/15/18	JLI	SW8260C
Acrolein	ND	4.8		ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	19		ug/Kg	1	12/15/18	JLI	SW8260C
Tert-butyl alcohol	ND	97		ug/Kg	1	12/15/18	JLI	SW8260C
<u>Semivolatiles</u>								
1,2,4,5-Tetrachlorobenzene	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
1,2,4-Trichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Diphenylhydrazine	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
1,3-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,4-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2,4,5-Trichlorophenol	ND	260	210	ug/Kg	1	12/14/18	WB	SW8270D
2,4,6-Trichlorophenol	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
2,4-Dichlorophenol	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dimethylphenol	ND	260	94	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrophenol	ND	260	260	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrotoluene	ND	190	150	ug/Kg	1	12/14/18	WB	SW8270D
2,6-Dinitrotoluene	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Chloronaphthalene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Chlorophenol	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylnaphthalene	250	J 260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylphenol (o-cresol)	ND	260	180	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitroaniline	ND	260	260	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitrophenol	ND	260	240	ug/Kg	1	12/14/18	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	260	150	ug/Kg	1	12/14/18	WB	SW8270D
3,3'-Dichlorobenzidine	ND	190	180	ug/Kg	1	12/14/18	WB	SW8270D
3-Nitroaniline	ND	380	760	ug/Kg	1	12/14/18	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	230	76	ug/Kg	1	12/14/18	WB	SW8270D
4-Bromophenyl phenyl ether	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloro-3-methylphenol	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloroaniline	ND	300	180	ug/Kg	1	12/14/18	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitroaniline	ND	380	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitrophenol	ND	380	170	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthene	490	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthylene	550	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Acetophenone	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Aniline	ND	300	300	ug/Kg	1	12/14/18	WB	SW8270D
Anthracene	1500	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Benz(a)anthracene	3400	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzidine	ND	380	220	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(a)pyrene	3200	190	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(b)fluoranthene	3000	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(ghi)perylene	1800	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(k)fluoranthene	2700	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzoic acid	ND	1900	760	ug/Kg	1	12/14/18	WB	SW8270D
Benzyl butyl phthalate	ND	260	97	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	260	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethyl)ether	ND	190	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Carbazole	360	190	150	ug/Kg	1	12/14/18	WB	SW8270D
Chrysene	3400	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenz(a,h)anthracene	720	190	120	ug/Kg	1	12/14/18	WB	SW8270D
Dibenzofuran	320	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Diethyl phthalate	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Dimethylphthalate	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-butylphthalate	ND	260	100	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-octylphthalate	ND	260	97	ug/Kg	1	12/14/18	WB	SW8270D
Fluoranthene	7000	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Fluorene	470	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobenzene	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Hexachlorobutadiene	ND	260	140	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorocyclopentadiene	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachloroethane	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Indeno(1,2,3-cd)pyrene	2100	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Isophorone	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Naphthalene	210	J 260	110	ug/Kg	1	12/14/18	WB	SW8270D
Nitrobenzene	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodimethylamine	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodiphenylamine	ND	260	150	ug/Kg	1	12/14/18	WB	SW8270D
Pentachloronitrobenzene	ND	260	140	ug/Kg	1	12/14/18	WB	SW8270D
Pentachlorophenol	ND	230	140	ug/Kg	1	12/14/18	WB	SW8270D
Phenanthrene	5000	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Phenol	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyrene	6500	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Pyridine	ND	260	93	ug/Kg	1	12/14/18	WB	SW8270D
<u>QA/QC Surrogates</u>								
% 2,4,6-Tribromophenol	68			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorobiphenyl	59			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorophenol	50			%	1	12/14/18	WB	30 - 130 %
% Nitrobenzene-d5	58			%	1	12/14/18	WB	30 - 130 %
% Phenol-d5	58			%	1	12/14/18	WB	30 - 130 %
% Terphenyl-d14	60			%	1	12/14/18	WB	30 - 130 %
Field Extraction	Completed					12/12/18		SW5035A

Parameter	Result	RL/ PQL	LOD/ MDL	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.

B = Present in blank, no bias suspected.

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit
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:

Per 1.4.6 of EPA method 8270D, 1,2-Diphenylhydrazine is unstable and readily converts to Azobenzene. Azobenzene is used for the calibration of 1,2-Diphenylhydrazine.

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

Volatile Comment:

There was a suppression of the last internal standard in the low level analysis, all affected compounds are reported from the methanol preserved high level analysis which did not exhibit this interference.

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

S - Laboratory solvent, contamination is possible.

If there are any questions regarding this data, please call Phoenix Client Services.
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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

11:45
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13977

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: SB 2 (7-9')

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.40	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Aluminum	9320	40		mg/Kg	10	12/15/18	CPP	SW6010C
Arsenic	8.39	0.80		mg/Kg	1	12/15/18	CPP	SW6010C
Barium	285	0.8		mg/Kg	1	12/15/18	EK	SW6010C
Beryllium	0.52	0.32		mg/Kg	1	12/15/18	CPP	SW6010C
Calcium	44500	40		mg/Kg	10	12/15/18	CPP	SW6010C
Cadmium	2.61	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Cobalt	9.08	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Chromium	23.2	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Copper	159	0.8		mg/kg	1	12/15/18	CPP	SW6010C
Iron	39300	40		mg/Kg	10	12/15/18	CPP	SW6010C
Mercury	13.0	1.4		mg/Kg	1	12/14/18	RS	SW7471B
Potassium	1840	8		mg/Kg	1	12/15/18	EK	SW6010C
Magnesium	4420	4.0		mg/Kg	1	12/15/18	CPP	SW6010C
Manganese	398	4.0		mg/Kg	10	12/15/18	CPP	SW6010C
Sodium	396	8		mg/Kg	1	12/15/18	CPP	SW6010C
Nickel	18.5	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Lead	959	8.0		mg/Kg	10	12/15/18	CPP	SW6010C
Antimony	< 4.0	4.0		mg/Kg	1	12/15/18	CPP	SW6010C
Selenium	< 1.6	1.6		mg/Kg	1	12/15/18	CPP	SW6010C
Thallium	< 1.6	1.6		mg/Kg	1	12/15/18	CPP	SW6010C
Vanadium	25.3	0.40		mg/Kg	1	12/15/18	CPP	SW6010C
Zinc	877	8.0		mg/Kg	10	12/15/18	CPP	SW6010C
Percent Solid	84			%		12/13/18	AK	SW846-%Solid
Soil Extraction for SVOA	Completed					12/13/18	JJ/CK	SW3545A
Mercury Digestion	Completed					12/14/18	EV/EV	SW7471B
Total Metals Digest	Completed					12/14/18	M/AG	SW3050B

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,1-Trichloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2,2-Tetrachloroethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2-Trichloroethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichlorobenzene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichloropropane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trichlorobenzene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trimethylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromo-3-chloropropane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromoethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichlorobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloropropane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,3,5-Trimethylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichlorobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichloropropane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
1,4-Dichlorobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
2,2-Dichloropropane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
2-Chlorotoluene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
2-Hexanone	ND	24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
2-Isopropyltoluene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
4-Chlorotoluene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
4-Methyl-2-pentanone	ND	24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Acetone	23	JS 24	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	9.6	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Benzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Bromobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Bromochloromethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Bromodichloromethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Bromoform	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Bromomethane	ND	4.8	1.9	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon Disulfide	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon tetrachloride	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Chlorobenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroform	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Chloromethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,2-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,3-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromochloromethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromomethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Dichlorodifluoromethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Ethylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Hexachlorobutadiene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Isopropylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
m&p-Xylene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl Ethyl Ketone	ND	29	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl t-butyl ether (MTBE)	ND	9.6	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Methylene chloride	ND	4.8	4.8	ug/Kg	1	12/15/18	JLI	SW8260C
Naphthalene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
n-Butylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
n-Propylbenzene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
o-Xylene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
p-Isopropyltoluene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
sec-Butylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Styrene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
tert-Butylbenzene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrachloroethene	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrahydrofuran (THF)	ND	9.6	2.4	ug/Kg	1	12/15/18	JLI	SW8260C
Toluene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,2-Dichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,3-Dichloropropene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,4-dichloro-2-butene	ND	9.6	2.4	ug/Kg	1	12/15/18	JLI	SW8260C
Trichloroethene	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorofluoromethane	ND	4.8	0.96	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorotrifluoroethane	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
Vinyl chloride	ND	4.8	0.48	ug/Kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	100			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	95			%	1	12/15/18	JLI	70 - 130 %
% Dibromofluoromethane	95			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	100			%	1	12/15/18	JLI	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	72		ug/kg	1	12/15/18	JLI	SW8260C
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	19		ug/Kg	1	12/15/18	JLI	SW8260C
Acrolein	ND	4.8		ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	19		ug/Kg	1	12/15/18	JLI	SW8260C
Tert-butyl alcohol	ND	96		ug/Kg	1	12/15/18	JLI	SW8260C
<u>Semivolatiles</u>								
1,2,4,5-Tetrachlorobenzene	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
1,2,4-Trichlorobenzene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Dichlorobenzene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Diphenylhydrazine	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
1,3-Dichlorobenzene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
1,4-Dichlorobenzene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
2,4,5-Trichlorophenol	ND	270	210	ug/Kg	1	12/14/18	WB	SW8270D
2,4,6-Trichlorophenol	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dichlorophenol	ND	190	140	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dimethylphenol	ND	270	96	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrophenol	ND	270	270	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrotoluene	ND	190	150	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
2,6-Dinitrotoluene	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Chloronaphthalene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Chlorophenol	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylnaphthalene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylphenol (o-cresol)	ND	270	180	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitroaniline	ND	270	270	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitrophenol	ND	270	250	ug/Kg	1	12/14/18	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	270	150	ug/Kg	1	12/14/18	WB	SW8270D
3,3'-Dichlorobenzidine	ND	190	180	ug/Kg	1	12/14/18	WB	SW8270D
3-Nitroaniline	ND	390	770	ug/Kg	1	12/14/18	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	230	77	ug/Kg	1	12/14/18	WB	SW8270D
4-Bromophenyl phenyl ether	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloro-3-methylphenol	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloroaniline	ND	310	180	ug/Kg	1	12/14/18	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitroaniline	ND	390	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitrophenol	ND	390	170	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthylene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Acetophenone	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Aniline	ND	310	310	ug/Kg	1	12/14/18	WB	SW8270D
Anthracene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benz(a)anthracene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzdine	ND	390	230	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(a)pyrene	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(b)fluoranthene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(ghi)perylene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(k)fluoranthene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzoic acid	ND	1900	770	ug/Kg	1	12/14/18	WB	SW8270D
Benzyl butyl phthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethyl)ether	ND	190	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Carbazole	ND	190	150	ug/Kg	1	12/14/18	WB	SW8270D
Chrysene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenz(a,h)anthracene	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenzofuran	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Diethyl phthalate	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Dimethylphthalate	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-butylphthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-octylphthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Fluoranthene	320	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Fluorene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobenzene	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobutadiene	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorocyclopentadiene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachloroethane	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
Indeno(1,2,3-cd)pyrene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D

1

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Isophorone	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Naphthalene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Nitrobenzene	ND	190	140	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodimethylamine	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodiphenylamine	ND	270	150	ug/Kg	1	12/14/18	WB	SW8270D
Pentachloronitrobenzene	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
Pentachlorophenol	ND	230	150	ug/Kg	1	12/14/18	WB	SW8270D
Phenanthrene	300	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Phenol	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyrene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Pyridine	ND	270	95	ug/Kg	1	12/14/18	WB	SW8270D
QA/QC Surrogates								
% 2,4,6-Tribromophenol	73			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorobiphenyl	63			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorophenol	50			%	1	12/14/18	WB	30 - 130 %
% Nitrobenzene-d5	60			%	1	12/14/18	WB	30 - 130 %
% Phenol-d5	61			%	1	12/14/18	WB	30 - 130 %
% Terphenyl-d14	65			%	1	12/14/18	WB	30 - 130 %
Field Extraction	Completed					12/12/18		SW5035A

1

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.

B = Present in blank, no bias suspected.

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit

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

Comments:

Per 1.4.6 of EPA method 8270D, 1,2-Diphenylhydrazine is unstable and readily converts to Azobenzene. Azobenzene is used for the calibration of 1,2-Diphenylhydrazine.

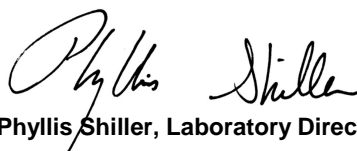
Please be advised that the NY 375 soil criteria for chromium are based on hexavalent chromium and trivalent chromium.

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

S - Laboratory solvent, contamination is possible.

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

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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

10:30
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13978

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: SB 3 (7-9')

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.37	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Aluminum	9190	37		mg/Kg	10	12/15/18	CPP	SW6010C
Arsenic	8.23	0.74		mg/Kg	1	12/15/18	CPP	SW6010C
Barium	203	0.7		mg/Kg	1	12/15/18	EK	SW6010C
Beryllium	0.56	0.29		mg/Kg	1	12/15/18	CPP	SW6010C
Calcium	6280	3.7		mg/Kg	1	12/15/18	CPP	SW6010C
Cadmium	0.90	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Cobalt	8.51	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Chromium	34.8	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Copper	63.9	0.7		mg/kg	1	12/15/18	CPP	SW6010C
Iron	25000	37		mg/Kg	10	12/15/18	CPP	SW6010C
Mercury	0.53	0.14		mg/Kg	1	12/14/18	RS	SW7471B
Potassium	1310	7		mg/Kg	1	12/15/18	EK	SW6010C
Magnesium	3000	3.7		mg/Kg	1	12/15/18	CPP	SW6010C
Manganese	228	3.7		mg/Kg	10	12/15/18	CPP	SW6010C
Sodium	375	7		mg/Kg	1	12/15/18	CPP	SW6010C
Nickel	43.2	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Lead	382	7.4		mg/Kg	10	12/15/18	CPP	SW6010C
Antimony	< 3.7	3.7		mg/Kg	1	12/15/18	CPP	SW6010C
Selenium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Thallium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Vanadium	23.5	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Zinc	195	7.4		mg/Kg	10	12/15/18	CPP	SW6010C
Percent Solid	87			%		12/13/18	AK	SW846-%Solid
Soil Extraction for SVOA	Completed					12/13/18	JJ/CK	SW3545A
Mercury Digestion	Completed					12/14/18	EV/EV	SW7471B
Total Metals Digest	Completed					12/14/18	M/AG	SW3050B

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,1-Trichloroethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2,2-Tetrachloroethane	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,1,2-Trichloroethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloropropene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichlorobenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,3-Trichloropropane	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,4-Trichlorobenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2,4-Trimethylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dibromo-3-chloropropane	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dibromoethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,2-Dichloroethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloropropane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
1,3,5-Trimethylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,3-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
1,3-Dichloropropane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
1,4-Dichlorobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
2,2-Dichloropropane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
2-Chlorotoluene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
2-Hexanone	ND	31	6.3	ug/Kg	1	12/15/18	JLI	SW8260C
2-Isopropyltoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
4-Chlorotoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
4-Methyl-2-pentanone	ND	31	6.3	ug/Kg	1	12/15/18	JLI	SW8260C
Acetone	23	JS 31	6.3	ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	13	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Benzene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Bromobenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Bromochloromethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Bromodichloromethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Bromoform	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Bromomethane	ND	6.3	2.5	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon Disulfide	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon tetrachloride	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Chlorobenzene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroform	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Chloromethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,2-Dichloroethene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,3-Dichloropropene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromochloromethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromomethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Dichlorodifluoromethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Ethylbenzene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Hexachlorobutadiene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Isopropylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
m&p-Xylene	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl Ethyl Ketone	ND	38	6.3	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl t-butyl ether (MTBE)	ND	13	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Methylene chloride	ND	6.3	6.3	ug/Kg	1	12/15/18	JLI	SW8260C
Naphthalene	71	J 320	64	ug/Kg	50	12/15/18	JLI	SW8260C
n-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
n-Propylbenzene	ND	320	64	ug/Kg	50	12/15/18	JLI	SW8260C
o-Xylene	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
p-Isopropyltoluene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
sec-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Styrene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
tert-Butylbenzene	ND	320	32	ug/Kg	50	12/15/18	JLI	SW8260C
Tetrachloroethene	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrahydrofuran (THF)	ND	13	3.1	ug/Kg	1	12/15/18	JLI	SW8260C
Toluene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,2-Dichloroethene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,3-Dichloropropene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,4-dichloro-2-butene	ND	640	160	ug/Kg	50	12/15/18	JLI	SW8260C
Trichloroethene	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorofluoromethane	ND	6.3	1.3	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorotrifluoroethane	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
Vinyl chloride	ND	6.3	0.63	ug/Kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	100			%	50	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	99			%	50	12/15/18	JLI	70 - 130 %
% Dibromofluoromethane	98			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	95			%	1	12/15/18	JLI	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	94		ug/kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	105			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	82			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	95			%	1	12/15/18	JLI	70 - 130 %
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	25		ug/Kg	1	12/15/18	JLI	SW8260C
Acrolein	ND	6.3		ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	25		ug/Kg	1	12/15/18	JLI	SW8260C
Tert-butyl alcohol	ND	130		ug/Kg	1	12/15/18	JLI	SW8260C
<u>Semivolatiles</u>								
1,2,4,5-Tetrachlorobenzene	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
1,2,4-Trichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Diphenylhydrazine	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
1,3-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
1,4-Dichlorobenzene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2,4,5-Trichlorophenol	ND	260	210	ug/Kg	1	12/14/18	WB	SW8270D
2,4,6-Trichlorophenol	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
2,4-Dichlorophenol	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dimethylphenol	ND	260	93	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrophenol	ND	260	260	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrotoluene	ND	190	150	ug/Kg	1	12/14/18	WB	SW8270D
2,6-Dinitrotoluene	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Chloronaphthalene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Chlorophenol	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylnaphthalene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylphenol (o-cresol)	ND	260	180	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitroaniline	ND	260	260	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitrophenol	ND	260	240	ug/Kg	1	12/14/18	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	260	150	ug/Kg	1	12/14/18	WB	SW8270D
3,3'-Dichlorobenzidine	ND	190	180	ug/Kg	1	12/14/18	WB	SW8270D
3-Nitroaniline	ND	380	750	ug/Kg	1	12/14/18	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	230	75	ug/Kg	1	12/14/18	WB	SW8270D
4-Bromophenyl phenyl ether	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloro-3-methylphenol	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloroaniline	ND	300	180	ug/Kg	1	12/14/18	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	260	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitroaniline	ND	380	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitrophenol	ND	380	170	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthylene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Acetophenone	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Aniline	ND	300	300	ug/Kg	1	12/14/18	WB	SW8270D
Anthracene	140	J 260	120	ug/Kg	1	12/14/18	WB	SW8270D
Benz(a)anthracene	430	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzidine	ND	380	220	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(a)pyrene	470	190	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(b)fluoranthene	400	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(ghi)perylene	330	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(k)fluoranthene	380	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzoic acid	ND	1900	750	ug/Kg	1	12/14/18	WB	SW8270D
Benzyl butyl phthalate	ND	260	97	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	260	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethyl)ether	ND	190	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	260	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Carbazole	ND	190	150	ug/Kg	1	12/14/18	WB	SW8270D
Chrysene	450	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenz(a,h)anthracene	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
Dibenzofuran	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Diethyl phthalate	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Dimethylphthalate	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-butylphthalate	ND	260	100	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-octylphthalate	ND	260	97	ug/Kg	1	12/14/18	WB	SW8270D
Fluoranthene	780	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Fluorene	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobenzene	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Hexachlorobutadiene	ND	260	140	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorocyclopentadiene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Hexachloroethane	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Indeno(1,2,3-cd)pyrene	330	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Isophorone	ND	190	110	ug/Kg	1	12/14/18	WB	SW8270D
Naphthalene	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Nitrobenzene	ND	190	130	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodimethylamine	ND	260	110	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	190	120	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodiphenylamine	ND	260	140	ug/Kg	1	12/14/18	WB	SW8270D
Pentachloronitrobenzene	ND	260	140	ug/Kg	1	12/14/18	WB	SW8270D
Pentachlorophenol	ND	230	140	ug/Kg	1	12/14/18	WB	SW8270D
Phenanthrene	520	260	110	ug/Kg	1	12/14/18	WB	SW8270D
Phenol	ND	260	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyrene	740	260	130	ug/Kg	1	12/14/18	WB	SW8270D
Pyridine	ND	260	92	ug/Kg	1	12/14/18	WB	SW8270D
<u>QA/QC Surrogates</u>								
% 2,4,6-Tribromophenol	79			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorobiphenyl	61			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorophenol	55			%	1	12/14/18	WB	30 - 130 %
% Nitrobenzene-d5	57			%	1	12/14/18	WB	30 - 130 %
% Phenol-d5	61			%	1	12/14/18	WB	30 - 130 %
% Terphenyl-d14	60			%	1	12/14/18	WB	30 - 130 %
Field Extraction	Completed					12/12/18		SW5035A

Parameter	Result	RL/ PQL	LOD/ MDL	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.

B = Present in blank, no bias suspected.

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit
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:

Per 1.4.6 of EPA method 8270D, 1,2-Diphenylhydrazine is unstable and readily converts to Azobenzene. Azobenzene is used for the calibration of 1,2-Diphenylhydrazine.

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

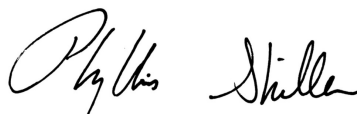
Volatile Comment:

There was a suppression of the last internal standard in the low level analysis, all affected compounds are reported from the methanol preserved high level analysis which did not exhibit this interference.

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

S - Laboratory solvent, contamination is possible.

If there are any questions regarding this data, please call Phoenix Client Services.
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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

9:45
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13979

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: SB 4 (8-10')

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.38	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Aluminum	9910	38		mg/Kg	10	12/15/18	CPP	SW6010C
Arsenic	5.21	0.76		mg/Kg	1	12/15/18	CPP	SW6010C
Barium	71.1	0.8		mg/Kg	1	12/15/18	EK	SW6010C
Beryllium	0.49	0.30		mg/Kg	1	12/15/18	CPP	SW6010C
Calcium	1540	3.8		mg/Kg	1	12/15/18	CPP	SW6010C
Cadmium	0.43	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Cobalt	10.7	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Chromium	17.4	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Copper	26.8	0.8		mg/kg	1	12/15/18	CPP	SW6010C
Iron	20000	38		mg/Kg	10	12/15/18	CPP	SW6010C
Mercury	0.24	0.13		mg/Kg	1	12/14/18	RS	SW7471B
Potassium	2030	8		mg/Kg	1	12/15/18	EK	SW6010C
Magnesium	2980	3.8		mg/Kg	1	12/15/18	CPP	SW6010C
Manganese	293	3.8		mg/Kg	10	12/15/18	CPP	SW6010C
Sodium	125	8		mg/Kg	1	12/15/18	CPP	SW6010C
Nickel	16.3	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Lead	62.6	0.8		mg/Kg	1	12/15/18	CPP	SW6010C
Antimony	< 3.8	3.8		mg/Kg	1	12/15/18	CPP	SW6010C
Selenium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Thallium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Vanadium	26.3	0.38		mg/Kg	1	12/15/18	CPP	SW6010C
Zinc	85.6	0.8		mg/Kg	1	12/15/18	CPP	SW6010C
Percent Solid	90			%		12/13/18	AK	SW846-%Solid
Soil Extraction for SVOA	Completed					12/13/18	JJ/CK	SW3545A
Mercury Digestion	Completed					12/14/18	EV/EV	SW7471B
Total Metals Digest	Completed					12/14/18	M/AG	SW3050B

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,1-Trichloroethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2,2-Tetrachloroethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2-Trichloroethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloropropene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichlorobenzene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichloropropane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trichlorobenzene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trimethylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromo-3-chloropropane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromoethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichlorobenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloroethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloropropane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,3,5-Trimethylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichlorobenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichloropropane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
1,4-Dichlorobenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
2,2-Dichloropropane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
2-Chlorotoluene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
2-Hexanone	ND	28	5.7	ug/Kg	1	12/15/18	JLI	SW8260C
2-Isopropyltoluene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
4-Chlorotoluene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
4-Methyl-2-pentanone	ND	28	5.7	ug/Kg	1	12/15/18	JLI	SW8260C
Acetone	10	JS 28	5.7	ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	11	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Benzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Bromobenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Bromochloromethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Bromodichloromethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Bromoform	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Bromomethane	ND	5.7	2.3	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon Disulfide	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon tetrachloride	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Chlorobenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroform	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Chloromethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,2-Dichloroethene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,3-Dichloropropene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromochloromethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromomethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Dichlorodifluoromethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Ethylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Hexachlorobutadiene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Isopropylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
m&p-Xylene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl Ethyl Ketone	ND	34	5.7	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl t-butyl ether (MTBE)	ND	11	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Methylene chloride	ND	5.7	5.7	ug/Kg	1	12/15/18	JLI	SW8260C
Naphthalene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
n-Butylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
n-Propylbenzene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
o-Xylene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
p-Isopropyltoluene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
sec-Butylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Styrene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
tert-Butylbenzene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrachloroethene	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrahydrofuran (THF)	ND	11	2.8	ug/Kg	1	12/15/18	JLI	SW8260C
Toluene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,2-Dichloroethene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,3-Dichloropropene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,4-dichloro-2-butene	ND	11	2.8	ug/Kg	1	12/15/18	JLI	SW8260C
Trichloroethene	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorofluoromethane	ND	5.7	1.1	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorotrifluoroethane	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
Vinyl chloride	ND	5.7	0.57	ug/Kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	112			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	90			%	1	12/15/18	JLI	70 - 130 %
% Dibromofluoromethane	94			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	98			%	1	12/15/18	JLI	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	85		ug/kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	112			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	90			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	98			%	1	12/15/18	JLI	70 - 130 %
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	23		ug/Kg	1	12/15/18	JLI	SW8260C
Acrolein	ND	5.7		ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	23		ug/Kg	1	12/15/18	JLI	SW8260C
Tert-butyl alcohol	ND	110		ug/Kg	1	12/15/18	JLI	SW8260C
<u>Semivolatiles</u>								
1,2,4,5-Tetrachlorobenzene	ND	250	130	ug/Kg	1	12/14/18	WB	SW8270D
1,2,4-Trichlorobenzene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Dichlorobenzene	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Diphenylhydrazine	ND	250	120	ug/Kg	1	12/14/18	WB	SW8270D
1,3-Dichlorobenzene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
1,4-Dichlorobenzene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
2,4,5-Trichlorophenol	ND	250	200	ug/Kg	1	12/14/18	WB	SW8270D
2,4,6-Trichlorophenol	ND	180	120	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
2,4-Dichlorophenol	ND	180	130	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dimethylphenol	ND	250	89	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrophenol	ND	250	250	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrotoluene	ND	180	140	ug/Kg	1	12/14/18	WB	SW8270D
2,6-Dinitrotoluene	ND	180	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Chloronaphthalene	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
2-Chlorophenol	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylnaphthalene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylphenol (o-cresol)	ND	250	170	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitroaniline	ND	250	250	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitrophenol	ND	250	230	ug/Kg	1	12/14/18	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	250	140	ug/Kg	1	12/14/18	WB	SW8270D
3,3'-Dichlorobenzidine	ND	180	170	ug/Kg	1	12/14/18	WB	SW8270D
3-Nitroaniline	ND	360	720	ug/Kg	1	12/14/18	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	220	72	ug/Kg	1	12/14/18	WB	SW8270D
4-Bromophenyl phenyl ether	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloro-3-methylphenol	ND	250	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloroaniline	ND	290	170	ug/Kg	1	12/14/18	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	250	120	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitroaniline	ND	360	120	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitrophenol	ND	360	160	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthylene	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
Acetophenone	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Aniline	ND	290	290	ug/Kg	1	12/14/18	WB	SW8270D
Anthracene	380	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Benz(a)anthracene	700	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzidine	ND	360	210	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(a)pyrene	640	180	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(b)fluoranthene	530	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(ghi)perylene	330	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(k)fluoranthene	500	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Benzoic acid	ND	1800	720	ug/Kg	1	12/14/18	WB	SW8270D
Benzyl butyl phthalate	ND	250	93	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	250	99	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethyl)ether	ND	180	97	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
Carbazole	ND	180	140	ug/Kg	1	12/14/18	WB	SW8270D
Chrysene	680	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Dibenz(a,h)anthracene	ND	180	120	ug/Kg	1	12/14/18	WB	SW8270D
Dibenzofuran	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Diethyl phthalate	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Dimethylphthalate	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-butylphthalate	ND	250	96	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-octylphthalate	ND	250	93	ug/Kg	1	12/14/18	WB	SW8270D
Fluoranthene	1700	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Fluorene	ND	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobenzene	ND	180	110	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Hexachlorobutadiene	ND	250	130	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorocyclopentadiene	ND	250	110	ug/Kg	1	12/14/18	WB	SW8270D
Hexachloroethane	ND	180	110	ug/Kg	1	12/14/18	WB	SW8270D
Indeno(1,2,3-cd)pyrene	360	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Isophorone	ND	180	100	ug/Kg	1	12/14/18	WB	SW8270D
Naphthalene	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
Nitrobenzene	ND	180	130	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodimethylamine	ND	250	100	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	180	120	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodiphenylamine	ND	250	140	ug/Kg	1	12/14/18	WB	SW8270D
Pentachloronitrobenzene	ND	250	130	ug/Kg	1	12/14/18	WB	SW8270D
Pentachlorophenol	ND	220	140	ug/Kg	1	12/14/18	WB	SW8270D
Phenanthrene	1300	250	100	ug/Kg	1	12/14/18	WB	SW8270D
Phenol	ND	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyrene	1500	250	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyridine	ND	250	89	ug/Kg	1	12/14/18	WB	SW8270D
QA/QC Surrogates								
% 2,4,6-Tribromophenol	74			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorobiphenyl	59			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorophenol	54			%	1	12/14/18	WB	30 - 130 %
% Nitrobenzene-d5	58			%	1	12/14/18	WB	30 - 130 %
% Phenol-d5	61			%	1	12/14/18	WB	30 - 130 %
% Terphenyl-d14	61			%	1	12/14/18	WB	30 - 130 %
Field Extraction	Completed					12/12/18		SW5035A

1

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.

B = Present in blank, no bias suspected.

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit

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:

Per 1.4.6 of EPA method 8270D, 1,2-Diphenylhydrazine is unstable and readily converts to Azobenzene. Azobenzene is used for the calibration of 1,2-Diphenylhydrazine.

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.

S - Laboratory solvent, contamination is possible.

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

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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

9:00
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13980

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: SB 5 (8-10')

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Silver	< 0.37	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Aluminum	12400	37		mg/Kg	10	12/15/18	CPP	SW6010C
Arsenic	2.64	0.74		mg/Kg	1	12/15/18	CPP	SW6010C
Barium	27.3	0.7		mg/Kg	1	12/15/18	EK	SW6010C
Beryllium	0.53	0.29		mg/Kg	1	12/15/18	CPP	SW6010C
Calcium	7670	3.7		mg/Kg	1	12/15/18	CPP	SW6010C
Cadmium	< 0.37	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Cobalt	6.10	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Chromium	17.9	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Copper	25.3	0.7		mg/kg	1	12/15/18	CPP	SW6010C
Iron	14700	37		mg/Kg	10	12/15/18	CPP	SW6010C
Mercury	0.03	0.03		mg/Kg	1	12/14/18	RS	SW7471B
Potassium	1470	7		mg/Kg	1	12/15/18	EK	SW6010C
Magnesium	6300	37		mg/Kg	10	12/15/18	CPP	SW6010C
Manganese	110	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Sodium	289	7		mg/Kg	1	12/15/18	CPP	SW6010C
Nickel	14.4	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Lead	11.7	0.7		mg/Kg	1	12/15/18	CPP	SW6010C
Antimony	< 3.7	3.7		mg/Kg	1	12/15/18	CPP	SW6010C
Selenium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Thallium	< 1.5	1.5		mg/Kg	1	12/15/18	CPP	SW6010C
Vanadium	22.5	0.37		mg/Kg	1	12/15/18	CPP	SW6010C
Zinc	64.4	0.7		mg/Kg	1	12/15/18	CPP	SW6010C
Percent Solid	85			%		12/13/18	AK	SW846-%Solid
Soil Extraction for SVOA	Completed					12/13/18	JJ/CK	SW3545A
Mercury Digestion	Completed					12/14/18	EV/EV	SW7471B
Total Metals Digest	Completed					12/14/18	M/AG	SW3050B

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,1-Trichloroethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2,2-Tetrachloroethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,1,2-Trichloroethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloroethene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,1-Dichloropropene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichlorobenzene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,3-Trichloropropane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trichlorobenzene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,2,4-Trimethylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dibromoethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichlorobenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloroethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,2-Dichloropropane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,3,5-Trimethylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichlorobenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
1,3-Dichloropropane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
1,4-Dichlorobenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
2,2-Dichloropropane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
2-Chlorotoluene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
2-Hexanone	ND	25	5.0	ug/Kg	1	12/15/18	JLI	SW8260C
2-Isopropyltoluene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
4-Chlorotoluene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
4-Methyl-2-pentanone	ND	25	5.0	ug/Kg	1	12/15/18	JLI	SW8260C
Acetone	36	S 25	5.0	ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	10	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Benzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Bromobenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Bromochloromethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Bromodichloromethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Bromoform	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Bromomethane	ND	5.0	2.0	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon Disulfide	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Carbon tetrachloride	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Chlorobenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Chloroform	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Chloromethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,2-Dichloroethene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
cis-1,3-Dichloropropene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromochloromethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Dibromomethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Dichlorodifluoromethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Ethylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Hexachlorobutadiene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Isopropylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
m&p-Xylene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl Ethyl Ketone	ND	30	5.0	ug/Kg	1	12/15/18	JLI	SW8260C
Methyl t-butyl ether (MTBE)	ND	10	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Methylene chloride	ND	5.0	5.0	ug/Kg	1	12/15/18	JLI	SW8260C
Naphthalene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
n-Butylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
n-Propylbenzene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
o-Xylene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
p-Isopropyltoluene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
sec-Butylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Styrene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
tert-Butylbenzene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrachloroethene	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Tetrahydrofuran (THF)	ND	10	2.5	ug/Kg	1	12/15/18	JLI	SW8260C
Toluene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,2-Dichloroethene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,3-Dichloropropene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
trans-1,4-dichloro-2-butene	ND	10	2.5	ug/Kg	1	12/15/18	JLI	SW8260C
Trichloroethene	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorofluoromethane	ND	5.0	1.0	ug/Kg	1	12/15/18	JLI	SW8260C
Trichlorotrifluoroethane	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
Vinyl chloride	ND	5.0	0.50	ug/Kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	101			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	96			%	1	12/15/18	JLI	70 - 130 %
% Dibromofluoromethane	94			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	100			%	1	12/15/18	JLI	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	75		ug/kg	1	12/15/18	JLI	SW8260C
<u>QA/QC Surrogates</u>								
% 1,2-dichlorobenzene-d4	101			%	1	12/15/18	JLI	70 - 130 %
% Bromofluorobenzene	96			%	1	12/15/18	JLI	70 - 130 %
% Toluene-d8	100			%	1	12/15/18	JLI	70 - 130 %
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	20		ug/Kg	1	12/15/18	JLI	SW8260C
Acrolein	ND	5.0		ug/Kg	1	12/15/18	JLI	SW8260C
Acrylonitrile	ND	20		ug/Kg	1	12/15/18	JLI	SW8260C
Tert-butyl alcohol	ND	100		ug/Kg	1	12/15/18	JLI	SW8260C
<u>Semivolatiles</u>								
1,2,4,5-Tetrachlorobenzene	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
1,2,4-Trichlorobenzene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Dichlorobenzene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
1,2-Diphenylhydrazine	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
1,3-Dichlorobenzene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
1,4-Dichlorobenzene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
2,4,5-Trichlorophenol	ND	270	210	ug/Kg	1	12/14/18	WB	SW8270D
2,4,6-Trichlorophenol	ND	200	120	ug/Kg	1	12/14/18	WB	SW8270D

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
2,4-Dichlorophenol	ND	200	140	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dimethylphenol	ND	270	97	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrophenol	ND	270	270	ug/Kg	1	12/14/18	WB	SW8270D
2,4-Dinitrotoluene	ND	200	150	ug/Kg	1	12/14/18	WB	SW8270D
2,6-Dinitrotoluene	ND	200	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Chloronaphthalene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Chlorophenol	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylnaphthalene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
2-Methylphenol (o-cresol)	ND	270	180	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitroaniline	ND	270	270	ug/Kg	1	12/14/18	WB	SW8270D
2-Nitrophenol	ND	270	250	ug/Kg	1	12/14/18	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	270	150	ug/Kg	1	12/14/18	WB	SW8270D
3,3'-Dichlorobenzidine	ND	200	180	ug/Kg	1	12/14/18	WB	SW8270D
3-Nitroaniline	ND	390	780	ug/Kg	1	12/14/18	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	230	78	ug/Kg	1	12/14/18	WB	SW8270D
4-Bromophenyl phenyl ether	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloro-3-methylphenol	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
4-Chloroaniline	ND	310	180	ug/Kg	1	12/14/18	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitroaniline	ND	390	130	ug/Kg	1	12/14/18	WB	SW8270D
4-Nitrophenol	ND	390	180	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Acenaphthylene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Acetophenone	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Aniline	ND	310	310	ug/Kg	1	12/14/18	WB	SW8270D
Anthracene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benz(a)anthracene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzidine	ND	390	230	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(a)pyrene	ND	200	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(b)fluoranthene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(ghi)perylene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzo(k)fluoranthene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Benzoic acid	ND	2000	780	ug/Kg	1	12/14/18	WB	SW8270D
Benzyl butyl phthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroethyl)ether	ND	200	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Carbazole	ND	200	160	ug/Kg	1	12/14/18	WB	SW8270D
Chrysene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenz(a,h)anthracene	ND	200	130	ug/Kg	1	12/14/18	WB	SW8270D
Dibenzofuran	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Diethyl phthalate	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Dimethylphthalate	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-butylphthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Di-n-octylphthalate	ND	270	100	ug/Kg	1	12/14/18	WB	SW8270D
Fluoranthene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Fluorene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorobenzene	ND	200	110	ug/Kg	1	12/14/18	WB	SW8270D

1

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Hexachlorobutadiene	ND	270	140	ug/Kg	1	12/14/18	WB	SW8270D
Hexachlorocyclopentadiene	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Hexachloroethane	ND	200	120	ug/Kg	1	12/14/18	WB	SW8270D
Indeno(1,2,3-cd)pyrene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Isophorone	ND	200	110	ug/Kg	1	12/14/18	WB	SW8270D
Naphthalene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Nitrobenzene	ND	200	140	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodimethylamine	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	200	130	ug/Kg	1	12/14/18	WB	SW8270D
N-Nitrosodiphenylamine	ND	270	150	ug/Kg	1	12/14/18	WB	SW8270D
Pentachloronitrobenzene	ND	270	150	ug/Kg	1	12/14/18	WB	SW8270D
Pentachlorophenol	ND	230	150	ug/Kg	1	12/14/18	WB	SW8270D
Phenanthrene	ND	270	110	ug/Kg	1	12/14/18	WB	SW8270D
Phenol	ND	270	120	ug/Kg	1	12/14/18	WB	SW8270D
Pyrene	ND	270	130	ug/Kg	1	12/14/18	WB	SW8270D
Pyridine	ND	270	96	ug/Kg	1	12/14/18	WB	SW8270D
QA/QC Surrogates								
% 2,4,6-Tribromophenol	77			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorobiphenyl	62			%	1	12/14/18	WB	30 - 130 %
% 2-Fluorophenol	62			%	1	12/14/18	WB	30 - 130 %
% Nitrobenzene-d5	61			%	1	12/14/18	WB	30 - 130 %
% Phenol-d5	68			%	1	12/14/18	WB	30 - 130 %
% Terphenyl-d14	64			%	1	12/14/18	WB	30 - 130 %
Field Extraction	Completed					12/12/18		SW5035A

1

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.

B = Present in blank, no bias suspected.

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 LOD=Limit of Detection MDL=Method Detection Limit1

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:

Per 1.4.6 of EPA method 8270D, 1,2-Diphenylhydrazine is unstable and readily converts to Azobenzene. Azobenzene is used for the calibration of 1,2-Diphenylhydrazine.

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.

S - Laboratory solvent, contamination is possible.

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

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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: GROUND WATER
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

8:40
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13981

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: GW1

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,1-Trichloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloropropene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,3-Trichloropropane	ND	0.25	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,4-Trimethylbenzene	1.8	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	0.50	0.50	ug/L	1	12/15/18	MH	SW8260C
1,2-Dibromoethane	ND	0.25	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichloroethane	ND	0.60	0.50	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3,5-Trimethylbenzene	0.34	J 1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2-Hexanone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C
2-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
4-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
4-Methyl-2-pentanone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Acetone	3.9	JS 5.0	2.5	ug/L	1	12/15/18	MH	SW8260C
Acrolein	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C
Acrylonitrile	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C
Benzene	ND	0.70	0.25	ug/L	1	12/15/18	MH	SW8260C
Bromobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Bromochloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Bromodichloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Bromoform	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Bromomethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Carbon Disulfide	0.40	J 1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Carbon tetrachloride	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Chlorobenzene	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Chloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Chloroform	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Chloromethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
cis-1,2-Dichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/15/18	MH	SW8260C
Dibromochloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Dibromomethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Dichlorodifluoromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Ethylbenzene	0.32	J 1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Hexachlorobutadiene	ND	0.50	0.20	ug/L	1	12/15/18	MH	SW8260C
Isopropylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
m&p-Xylene	1.2	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Methyl ethyl ketone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Methylene chloride	ND	3.0	1.0	ug/L	1	12/15/18	MH	SW8260C
Naphthalene	1.6	1.0	1.0	ug/L	1	12/15/18	MH	SW8260C
n-Butylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
n-Propylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
o-Xylene	0.60	J 1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
p-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
sec-Butylbenzene	1.5	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Styrene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
tert-Butylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Tetrachloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Tetrahydrofuran (THF)	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C
Toluene	1.3	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
trans-1,2-Dichloroethene	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/15/18	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C
Trichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Trichlorofluoromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Trichlorotrifluoroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
Vinyl chloride	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
QA/QC Surrogates								
% 1,2-dichlorobenzene-d4	101			%	1	12/15/18	MH	70 - 130 %
% Bromofluorobenzene	96			%	1	12/15/18	MH	70 - 130 %
% Dibromofluoromethane	100			%	1	12/15/18	MH	70 - 130 %

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
% Toluene-d8	98			%	1	12/15/18	MH	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	100		ug/l	1	12/15/18	MH	SW8260C
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0		ug/L	1	12/15/18	MH	SW8260C
Acrolein	ND	5.0		ug/L	1	12/15/18	MH	SW8260C
Acrylonitrile	ND	5.0		ug/L	1	12/15/18	MH	SW8260C
Tert-butyl alcohol	ND	50		ug/L	1	12/15/18	MH	SW8260C

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit
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.

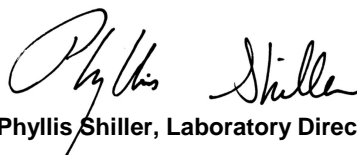
Comments:

Volatile Comment:

Where the LOD justifies lowering the RL/PQL, the RL/PQL of some compounds are evaluated below the lowest calibration standard in order to meet criteria.

S - Laboratory solvent, contamination is possible.

If there are any questions regarding this data, please call Phoenix Client Services.
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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date Time
 12/12/18 10:00
 12/13/18 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13982

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: GW2

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1,1-Trichloroethane	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1-Dichloroethane	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1-Dichloroethene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,1-Dichloropropene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2,3-Trichloropropane	ND	0.25	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	0.50	0.50	ug/L	1	12/14/18	MH	SW8260C
1,2-Dibromoethane	ND	0.25	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,2-Dichloroethane	ND	0.60	0.50	ug/L	1	12/14/18	MH	SW8260C
1,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,3-Dichloropropane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
2,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
2-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
2-Hexanone	ND	2.5	2.5	ug/L	1	12/14/18	MH	SW8260C
2-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
4-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
4-Methyl-2-pentanone	ND	2.5	2.5	ug/L	1	12/14/18	MH	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
Acetone	ND	5.0	2.5	ug/L	1	12/14/18	MH	SW8260C
Acrolein	ND	5.0	2.5	ug/L	1	12/14/18	MH	SW8260C
Acrylonitrile	ND	5.0	2.5	ug/L	1	12/14/18	MH	SW8260C
Benzene	ND	0.70	0.25	ug/L	1	12/14/18	MH	SW8260C
Bromobenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Bromochloromethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Bromodichloromethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Bromoform	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Bromomethane	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Carbon Disulfide	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Carbon tetrachloride	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Chlorobenzene	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Chloroethane	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Chloroform	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Chloromethane	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
cis-1,2-Dichloroethene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/14/18	MH	SW8260C
Dibromochloromethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Dibromomethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Dichlorodifluoromethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Ethylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Hexachlorobutadiene	ND	0.50	0.20	ug/L	1	12/14/18	MH	SW8260C
Isopropylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
m&p-Xylene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Methyl ethyl ketone	ND	2.5	2.5	ug/L	1	12/14/18	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Methylene chloride	ND	3.0	1.0	ug/L	1	12/14/18	MH	SW8260C
Naphthalene	ND	1.0	1.0	ug/L	1	12/14/18	MH	SW8260C
n-Butylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
n-Propylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
o-Xylene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
p-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
sec-Butylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Styrene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
tert-Butylbenzene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Tetrachloroethene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Tetrahydrofuran (THF)	ND	5.0	2.5	ug/L	1	12/14/18	MH	SW8260C
Toluene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
trans-1,2-Dichloroethene	ND	5.0	0.25	ug/L	1	12/14/18	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/14/18	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	2.5	2.5	ug/L	1	12/14/18	MH	SW8260C
Trichloroethene	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Trichlorofluoromethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Trichlorotrifluoroethane	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
Vinyl chloride	ND	1.0	0.25	ug/L	1	12/14/18	MH	SW8260C
QA/QC Surrogates								
% 1,2-dichlorobenzene-d4	99			%	1	12/14/18	MH	70 - 130 %
% Bromofluorobenzene	97			%	1	12/14/18	MH	70 - 130 %
% Dibromofluoromethane	98			%	1	12/14/18	MH	70 - 130 %

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
% Toluene-d8	99			%	1	12/14/18	MH	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	100		ug/l	1	12/14/18	MH	SW8260C
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0		ug/L	1	12/14/18	MH	SW8260C
Acrolein	ND	5.0		ug/L	1	12/14/18	MH	SW8260C
Acrylonitrile	ND	5.0		ug/L	1	12/14/18	MH	SW8260C
Tert-butyl alcohol	ND	50		ug/L	1	12/14/18	MH	SW8260C

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 LOD=Limit of Detection MDL=Method Detection Limit

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.

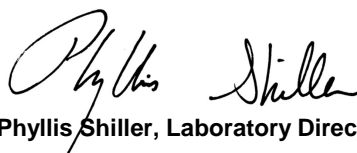
Comments:

Volatile Comment:

Where the LOD justifies lowering the RL/PQL, the RL/PQL of some compounds are evaluated below the lowest calibration standard in order to meet criteria.

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

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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab 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
 December 19, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: SOIL
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date

12/12/18
 12/13/18

Time

10:45
 16:30

Laboratory Data

SDG ID: GCC13976
 Phoenix ID: CC13983

Project ID: 188 EAST 135TH ST BRONX NY
 Client ID: GW3

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,1-Trichloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,1-Dichloropropene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,3-Trichloropropane	ND	0.25	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	0.50	0.50	ug/L	1	12/15/18	MH	SW8260C
1,2-Dibromoethane	ND	0.25	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichloroethane	ND	0.60	0.50	ug/L	1	12/15/18	MH	SW8260C
1,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,3-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2,2-Dichloropropane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
2-Hexanone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C
2-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
4-Chlorotoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C
4-Methyl-2-pentanone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference	
Acetone	3.7	JS 5.0	2.5	ug/L	1	12/15/18	MH	SW8260C	
Acrolein	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C	
Acrylonitrile	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C	
Benzene	ND	0.70	0.25	ug/L	1	12/15/18	MH	SW8260C	
Bromobenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Bromochloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Bromodichloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Bromoform	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Bromomethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Carbon Disulfide	0.39	J 1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Carbon tetrachloride	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Chlorobenzene	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Chloroethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Chloroform	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Chloromethane	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
cis-1,2-Dichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
cis-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/15/18	MH	SW8260C	
Dibromochloromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Dibromomethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Dichlorodifluoromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Ethylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Hexachlorobutadiene	ND	0.50	0.20	ug/L	1	12/15/18	MH	SW8260C	
Isopropylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
m&p-Xylene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Methyl ethyl ketone	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C	
Methyl t-butyl ether (MTBE)	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Methylene chloride	ND	3.0	1.0	ug/L	1	12/15/18	MH	SW8260C	
Naphthalene	ND	1.0	1.0	ug/L	1	12/15/18	MH	SW8260C	
n-Butylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
n-Propylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
o-Xylene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
p-Isopropyltoluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
sec-Butylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Styrene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
tert-Butylbenzene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Tetrachloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Tetrahydrofuran (THF)	ND	5.0	2.5	ug/L	1	12/15/18	MH	SW8260C	
Toluene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
trans-1,2-Dichloroethene	ND	5.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
trans-1,3-Dichloropropene	ND	0.40	0.25	ug/L	1	12/15/18	MH	SW8260C	
trans-1,4-dichloro-2-butene	ND	2.5	2.5	ug/L	1	12/15/18	MH	SW8260C	
Trichloroethene	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Trichlorofluoromethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Trichlorotrifluoroethane	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
Vinyl chloride	ND	1.0	0.25	ug/L	1	12/15/18	MH	SW8260C	
QA/QC Surrogates									
% 1,2-dichlorobenzene-d4	100			%	1	12/15/18	MH	70 - 130 %	
% Bromofluorobenzene	96			%	1	12/15/18	MH	70 - 130 %	
% Dibromofluoromethane	100			%	1	12/15/18	MH	70 - 130 %	

Parameter	Result	RL/ PQL	LOD/ MDL	Units	Dilution	Date/Time	By	Reference
% Toluene-d8	99			%	1	12/15/18	MH	70 - 130 %
<u>1,4-dioxane</u>								
1,4-dioxane	ND	100		ug/l	1	12/15/18	MH	SW8260C
<u>Volatiles</u>								
1,1,1,2-Tetrachloroethane	ND	1.0		ug/L	1	12/15/18	MH	SW8260C
Acrolein	ND	5.0		ug/L	1	12/15/18	MH	SW8260C
Acrylonitrile	ND	5.0		ug/L	1	12/15/18	MH	SW8260C
Tert-butyl alcohol	ND	50		ug/L	1	12/15/18	MH	SW8260C

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 J=Estimated Below RL LOD=Limit of Detection MDL=Method Detection Limit
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.

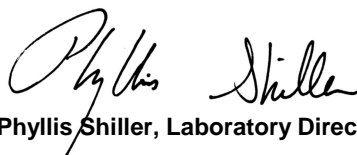
Comments:

Volatile Comment:

Where the LOD justifies lowering the RL/PQL, the RL/PQL of some compounds are evaluated below the lowest calibration standard in order to meet criteria.

S - Laboratory solvent, contamination is possible.

If there are any questions regarding this data, please call Phoenix Client Services.
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Phyllis Shiller, Laboratory Director

December 19, 2018

Reviewed and Released by: Greg Lawrence, Assistant Lab Director



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



QA/QC Report

December 19, 2018

QA/QC Data

SDG I.D.: GCC13976

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 459938 (mg/kg), QC Sample No: CC12609 (CC13976, CC13977, CC13978, CC13979, CC13980)

Mercury - Soil	BRL	0.02	0.20	0.29	NC	82.0	91.7	11.2	59.3	61.2	3.2	70 - 130	30 m
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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 460033 (mg/kg), QC Sample No: CC14009 (CC13976, CC13977, CC13978, CC13979, CC13980)

ICP Metals - Soil

Aluminum	BRL	5.1	9510	9690	1.90	110						NC	75 - 125	30
Antimony	BRL	3.4	<3.5	<3.4	NC	123						88.6	75 - 125	30
Arsenic	BRL	0.68	7.18	7.57	5.30	101						99.8	75 - 125	30
Barium	0.90	0.34	50.8	62.5	20.7	99.1						110	75 - 125	30
Beryllium	BRL	0.27	0.44	0.47	NC	96.7						101	75 - 125	30
Cadmium	BRL	0.34	<0.35	0.34	NC	98.3						101	75 - 125	30
Calcium	BRL	5.1	1510	1590	5.20	97.8						NC	75 - 125	30
Chromium	BRL	0.34	24.7	29.2	16.7	102						108	75 - 125	30
Cobalt	BRL	0.34	2.98	3.08	3.30	103						103	75 - 125	30
Copper	BRL	0.68	4.8	5.01	4.30	95.5						102	75 - 125	30
Iron	BRL	5.1	20000	20500	2.50	125						NC	75 - 125	30
Lead	BRL	0.34	5.6	5.82	3.90	101						102	75 - 125	30
Magnesium	BRL	5.1	1480	1570	5.90	106						NC	75 - 125	30
Manganese	BRL	0.34	62.7	63.9	1.90	101						107	75 - 125	30
Nickel	BRL	0.34	4.75	5.02	5.50	97.3						103	75 - 125	30
Potassium	BRL	5.1	1460	1580	7.90	113						NC	75 - 125	30
Selenium	BRL	1.4	<1.4	<1.4	NC	99.7						99.9	75 - 125	30
Silver	BRL	0.34	<0.35	<0.34	NC	95.3						101	75 - 125	30
Sodium	BRL	5.1	51	123	82.8	99.0						111	75 - 125	30 r
Thallium	BRL	3.1	<1.4	<3.1	NC	98.8						102	75 - 125	30
Vanadium	BRL	0.34	28.2	29.6	4.80	114						106	75 - 125	30
Zinc	BRL	0.68	23.0	25.5	10.3	97.8						101	75 - 125	30

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

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



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QA/QC Report

December 19, 2018

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
QA/QC Batch 459811 (ug/kg), QC Sample No: CC13734 (CC13976, CC13977, CC13978, CC13979, CC13980)										
Semivolatiles - Soil										
1,2,4,5-Tetrachlorobenzene	ND	230	73	70	4.2	73	69	5.6	30 - 130	30
1,2,4-Trichlorobenzene	ND	230	71	69	2.9	68	64	6.1	30 - 130	30
1,2-Dichlorobenzene	ND	180	65	63	3.1	61	58	5.0	30 - 130	30
1,2-Diphenylhydrazine	ND	230	68	66	3.0	70	66	5.9	30 - 130	30
1,3-Dichlorobenzene	ND	230	64	62	3.2	55	53	3.7	30 - 130	30
1,4-Dichlorobenzene	ND	230	65	63	3.1	59	56	5.2	30 - 130	30
2,4,5-Trichlorophenol	ND	230	84	77	8.7	86	80	7.2	30 - 130	30
2,4,6-Trichlorophenol	ND	130	82	78	5.0	84	78	7.4	30 - 130	30
2,4-Dichlorophenol	ND	130	80	78	2.5	80	77	3.8	30 - 130	30
2,4-Dimethylphenol	ND	230	81	79	2.5	85	81	4.8	30 - 130	30
2,4-Dinitrophenol	ND	230	14	<10	NC	31	34	9.2	30 - 130	30
2,4-Dinitrotoluene	ND	130	79	78	1.3	83	81	2.4	30 - 130	30
2,6-Dinitrotoluene	ND	130	81	80	1.2	84	82	2.4	30 - 130	30
2-Chloronaphthalene	ND	230	75	72	4.1	76	69	9.7	30 - 130	30
2-Chlorophenol	ND	230	72	71	1.4	76	70	8.2	30 - 130	30
2-Methylnaphthalene	ND	230	72	70	2.8	73	69	5.6	30 - 130	30
2-Methylphenol (o-cresol)	ND	230	70	69	1.4	76	72	5.4	30 - 130	30
2-Nitroaniline	ND	330	103	96	7.0	101	101	0.0	30 - 130	30
2-Nitrophenol	ND	230	75	76	1.3	76	73	4.0	30 - 130	30
3&4-Methylphenol (m&p-cresol)	ND	230	76	75	1.3	83	79	4.9	30 - 130	30
3,3'-Dichlorobenzidine	ND	130	78	80	2.5	96	90	6.5	30 - 130	30
3-Nitroaniline	ND	330	93	90	3.3	97	96	1.0	30 - 130	30
4,6-Dinitro-2-methylphenol	ND	230	35	22	45.6	44	47	6.6	30 - 130	30
4-Bromophenyl phenyl ether	ND	230	82	79	3.7	82	73	11.6	30 - 130	30
4-Chloro-3-methylphenol	ND	230	81	79	2.5	82	82	0.0	30 - 130	30
4-Chloroaniline	ND	230	74	74	0.0	81	78	3.8	30 - 130	30
4-Chlorophenyl phenyl ether	ND	230	81	77	5.1	79	74	6.5	30 - 130	30
4-Nitroaniline	ND	230	79	77	2.6	83	82	1.2	30 - 130	30
4-Nitrophenol	ND	230	76	74	2.7	79	77	2.6	30 - 130	30
Acenaphthene	ND	230	76	72	5.4	74	70	5.6	30 - 130	30
Acenaphthylene	ND	130	70	68	2.9	71	66	7.3	30 - 130	30
Acetophenone	ND	230	62	61	1.6	70	65	7.4	30 - 130	30
Aniline	ND	330	58	58	0.0	110	60	58.8	30 - 130	30
Anthracene	ND	230	76	75	1.3	79	72	9.3	30 - 130	30
Benz(a)anthracene	ND	230	75	74	1.3	73	67	8.6	30 - 130	30
Benzidine	ND	330	55	55	0.0	45	61	30.2	30 - 130	30
Benzo(a)pyrene	ND	130	75	73	2.7	73	66	10.1	30 - 130	30
Benzo(b)fluoranthene	ND	160	79	77	2.6	72	68	5.7	30 - 130	30
Benzo(ghi)perylene	ND	230	72	71	1.4	76	68	11.1	30 - 130	30
Benzo(k)fluoranthene	ND	230	76	72	5.4	71	65	8.8	30 - 130	30
Benzoic Acid	ND	330	<10	<10	NC	35	35	0.0	30 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
Benzyl butyl phthalate	ND	230	84	81	3.6	81	73	10.4	30 - 130	30
Bis(2-chloroethoxy)methane	ND	230	71	71	0.0	70	66	5.9	30 - 130	30
Bis(2-chloroethyl)ether	ND	130	60	58	3.4	61	57	6.8	30 - 130	30
Bis(2-chloroisopropyl)ether	ND	230	53	52	1.9	52	49	5.9	30 - 130	30
Bis(2-ethylhexyl)phthalate	ND	230	86	83	3.6	90	78	14.3	30 - 130	30
Carbazole	ND	230	77	75	2.6	81	74	9.0	30 - 130	30
Chrysene	ND	230	78	75	3.9	78	71	9.4	30 - 130	30
Dibenz(a,h)anthracene	ND	130	80	79	1.3	85	75	12.5	30 - 130	30
Dibenzofuran	ND	230	76	72	5.4	77	72	6.7	30 - 130	30
Diethyl phthalate	ND	230	80	77	3.8	79	75	5.2	30 - 130	30
Dimethylphthalate	ND	230	80	77	3.8	80	75	6.5	30 - 130	30
Di-n-butylphthalate	ND	670	84	82	2.4	83	76	8.8	30 - 130	30
Di-n-octylphthalate	ND	230	87	84	3.5	86	78	9.8	30 - 130	30
Fluoranthene	ND	230	79	76	3.9	77	72	6.7	30 - 130	30
Fluorene	ND	230	78	74	5.3	77	73	5.3	30 - 130	30
Hexachlorobenzene	ND	130	75	73	2.7	73	64	13.1	30 - 130	30
Hexachlorobutadiene	ND	230	73	72	1.4	66	63	4.7	30 - 130	30
Hexachlorocyclopentadiene	ND	230	53	53	0.0	22	12	58.8	30 - 130	30
Hexachloroethane	ND	130	63	61	3.2	53	50	5.8	30 - 130	30
Indeno(1,2,3-cd)pyrene	ND	230	71	72	1.4	76	68	11.1	30 - 130	30
Isophorone	ND	130	66	65	1.5	65	62	4.7	30 - 130	30
Naphthalene	ND	230	69	68	1.5	69	65	6.0	30 - 130	30
Nitrobenzene	ND	130	68	68	0.0	72	67	7.2	30 - 130	30
N-Nitrosodimethylamine	ND	230	59	57	3.4	63	58	8.3	30 - 130	30
N-Nitrosodi-n-propylamine	ND	130	70	69	1.4	73	68	7.1	30 - 130	30
N-Nitrosodiphenylamine	ND	130	80	77	3.8	79	75	5.2	30 - 130	30
Pentachloronitrobenzene	ND	230	77	76	1.3	79	72	9.3	30 - 130	30
Pentachlorophenol	ND	230	78	71	9.4	86	81	6.0	30 - 130	30
Phenanthrene	ND	130	76	73	4.0	77	73	5.3	30 - 130	30
Phenol	ND	230	72	70	2.8	75	72	4.1	30 - 130	30
Pyrene	ND	230	79	77	2.6	79	74	6.5	30 - 130	30
Pyridine	ND	230	44	42	4.7	44	42	4.7	30 - 130	30
% 2,4,6-Tribromophenol	61	%	74	71	4.1	79	72	9.3	30 - 130	30
% 2-Fluorobiphenyl	58	%	62	61	1.6	64	59	8.1	30 - 130	30
% 2-Fluorophenol	55	%	66	64	3.1	72	65	10.2	30 - 130	30
% Nitrobenzene-d5	53	%	58	58	0.0	63	59	6.6	30 - 130	30
% Phenol-d5	61	%	70	68	2.9	74	71	4.1	30 - 130	30
% Terphenyl-d14	54	%	62	60	3.3	63	57	10.0	30 - 130	30

m,r

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 460180 (ug/L), QC Sample No: CC12929 (CC13981, CC13983)

Volatiles - Ground Water

1,1,1,2-Tetrachloroethane	ND	1.0	109	109	0.0				70 - 130	30
1,1,1-Trichloroethane	ND	1.0	100	100	0.0				70 - 130	30
1,1,2,2-Tetrachloroethane	ND	0.50	111	108	2.7				70 - 130	30
1,1,2-Trichloroethane	ND	1.0	107	110	2.8				70 - 130	30
1,1-Dichloroethane	ND	1.0	98	99	1.0				70 - 130	30
1,1-Dichloroethene	ND	1.0	100	101	1.0				70 - 130	30
1,1-Dichloropropene	ND	1.0	103	101	2.0				70 - 130	30
1,2,3-Trichlorobenzene	ND	1.0	115	115	0.0				70 - 130	30
1,2,3-Trichloropropane	ND	1.0	105	103	1.9				70 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
1,2,4-Trichlorobenzene	ND	1.0	108	110	1.8				70 - 130	30
1,2,4-Trimethylbenzene	ND	1.0	99	100	1.0				70 - 130	30
1,2-Dibromo-3-chloropropane	ND	1.0	118	124	5.0				70 - 130	30
1,2-Dibromoethane	ND	1.0	111	108	2.7				70 - 130	30
1,2-Dichlorobenzene	ND	1.0	105	106	0.9				70 - 130	30
1,2-Dichloroethane	ND	1.0	106	106	0.0				70 - 130	30
1,2-Dichloropropane	ND	1.0	103	101	2.0				70 - 130	30
1,3,5-Trimethylbenzene	ND	1.0	98	99	1.0				70 - 130	30
1,3-Dichlorobenzene	ND	1.0	101	102	1.0				70 - 130	30
1,3-Dichloropropane	ND	1.0	108	106	1.9				70 - 130	30
1,4-Dichlorobenzene	ND	1.0	102	103	1.0				70 - 130	30
1,4-dioxane	ND	100	109	110	0.9				70 - 130	30
2,2-Dichloropropane	ND	1.0	106	105	0.9				70 - 130	30
2-Chlorotoluene	ND	1.0	98	100	2.0				70 - 130	30
2-Hexanone	ND	5.0	87	86	1.2				70 - 130	30
2-Isopropyltoluene	ND	1.0	101	103	2.0				70 - 130	30
4-Chlorotoluene	ND	1.0	98	99	1.0				70 - 130	30
4-Methyl-2-pentanone	ND	5.0	95	93	2.1				70 - 130	30
Acetone	ND	5.0	75	75	0.0				70 - 130	30
Acrolein	ND	5.0	110	110	0.0				70 - 130	30
Acrylonitrile	ND	5.0	108	104	3.8				70 - 130	30
Benzene	ND	0.70	100	98	2.0				70 - 130	30
Bromobenzene	ND	1.0	102	101	1.0				70 - 130	30
Bromochloromethane	ND	1.0	104	102	1.9				70 - 130	30
Bromodichloromethane	ND	0.50	111	108	2.7				70 - 130	30
Bromoform	ND	1.0	119	112	6.1				70 - 130	30
Bromomethane	ND	1.0	79	81	2.5				70 - 130	30
Carbon Disulfide	ND	1.0	99	97	2.0				70 - 130	30
Carbon tetrachloride	ND	1.0	100	99	1.0				70 - 130	30
Chlorobenzene	ND	1.0	100	101	1.0				70 - 130	30
Chloroethane	ND	1.0	103	106	2.9				70 - 130	30
Chloroform	ND	1.0	99	96	3.1				70 - 130	30
Chloromethane	ND	1.0	84	84	0.0				70 - 130	30
cis-1,2-Dichloroethene	ND	1.0	101	102	1.0				70 - 130	30
cis-1,3-Dichloropropene	ND	0.40	109	106	2.8				70 - 130	30
Dibromochloromethane	ND	0.50	117	117	0.0				70 - 130	30
Dibromomethane	ND	1.0	108	104	3.8				70 - 130	30
Dichlorodifluoromethane	ND	1.0	104	105	1.0				70 - 130	30
Ethylbenzene	ND	1.0	101	102	1.0				70 - 130	30
Hexachlorobutadiene	ND	0.40	103	107	3.8				70 - 130	30
Isopropylbenzene	ND	1.0	98	99	1.0				70 - 130	30
m&p-Xylene	ND	1.0	100	100	0.0				70 - 130	30
Methyl ethyl ketone	ND	5.0	89	89	0.0				70 - 130	30
Methyl t-butyl ether (MTBE)	ND	1.0	109	112	2.7				70 - 130	30
Methylene chloride	ND	1.0	99	98	1.0				70 - 130	30
Naphthalene	ND	1.0	119	117	1.7				70 - 130	30
n-Butylbenzene	ND	1.0	101	103	2.0				70 - 130	30
n-Propylbenzene	ND	1.0	98	99	1.0				70 - 130	30
o-Xylene	ND	1.0	103	102	1.0				70 - 130	30
p-Isopropyltoluene	ND	1.0	99	101	2.0				70 - 130	30
sec-Butylbenzene	ND	1.0	102	104	1.9				70 - 130	30
Styrene	ND	1.0	103	103	0.0				70 - 130	30
tert-butyl alcohol	ND	10	113	121	6.8				70 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blk		LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
	Blank	RL								
tert-Butylbenzene	ND	1.0	97	99	2.0				70 - 130	30
Tetrachloroethene	ND	1.0	102	102	0.0				70 - 130	30
Tetrahydrofuran (THF)	ND	2.5	102	103	1.0				70 - 130	30
Toluene	ND	1.0	101	100	1.0				70 - 130	30
trans-1,2-Dichloroethene	ND	1.0	101	101	0.0				70 - 130	30
trans-1,3-Dichloropropene	ND	0.40	106	103	2.9				70 - 130	30
trans-1,4-dichloro-2-butene	ND	5.0	116	115	0.9				70 - 130	30
Trichloroethene	ND	1.0	103	102	1.0				70 - 130	30
Trichlorofluoromethane	ND	1.0	104	102	1.9				70 - 130	30
Trichlorotrifluoroethane	ND	1.0	99	104	4.9				70 - 130	30
Vinyl chloride	ND	1.0	102	100	2.0				70 - 130	30
% 1,2-dichlorobenzene-d4	103	%	102	101	1.0				70 - 130	30
% Bromofluorobenzene	97	%	102	102	0.0				70 - 130	30
% Dibromofluoromethane	100	%	101	100	1.0				70 - 130	30
% Toluene-d8	98	%	101	100	1.0				70 - 130	30

Comment:

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

Additional 8260 criteria: 10% of LCS/LCSD compounds can be outside of acceptance criteria as long as recovery is 40-160%.

QA/QC Batch 460184 (ug/L), QC Sample No: CC13632 (CC13982)

Volatiles

1,1,1,2-Tetrachloroethane	ND	1.0	96	104	8.0				70 - 130	30
1,1,1-Trichloroethane	ND	1.0	87	97	10.9				70 - 130	30
1,1,2,2-Tetrachloroethane	ND	0.50	93	105	12.1				70 - 130	30
1,1,2-Trichloroethane	ND	1.0	93	105	12.1				70 - 130	30
1,1-Dichloroethane	ND	1.0	88	96	8.7				70 - 130	30
1,1-Dichloroethene	ND	1.0	91	99	8.4				70 - 130	30
1,1-Dichloropropene	ND	1.0	90	98	8.5				70 - 130	30
1,2,3-Trichlorobenzene	ND	1.0	93	106	13.1				70 - 130	30
1,2,3-Trichloropropane	ND	1.0	89	96	7.6				70 - 130	30
1,2,4-Trichlorobenzene	ND	1.0	92	103	11.3				70 - 130	30
1,2,4-Trimethylbenzene	ND	1.0	91	95	4.3				70 - 130	30
1,2-Dibromo-3-chloropropane	ND	1.0	98	107	8.8				70 - 130	30
1,2-Dibromoethane	ND	1.0	93	104	11.2				70 - 130	30
1,2-Dichlorobenzene	ND	1.0	93	100	7.3				70 - 130	30
1,2-Dichloroethane	ND	1.0	87	101	14.9				70 - 130	30
1,2-Dichloropropane	ND	1.0	88	97	9.7				70 - 130	30
1,3,5-Trimethylbenzene	ND	1.0	91	95	4.3				70 - 130	30
1,3-Dichlorobenzene	ND	1.0	92	98	6.3				70 - 130	30
1,3-Dichloropropane	ND	1.0	92	102	10.3				70 - 130	30
1,4-Dichlorobenzene	ND	1.0	91	98	7.4				70 - 130	30
1,4-dioxane	ND	100	96	94	2.1				70 - 130	30
2,2-Dichloropropane	ND	1.0	90	96	6.5				70 - 130	30
2-Chlorotoluene	ND	1.0	92	96	4.3				70 - 130	30
2-Hexanone	ND	5.0	71	80	11.9				70 - 130	30
2-Isopropyltoluene	ND	1.0	94	98	4.2				70 - 130	30
4-Chlorotoluene	ND	1.0	90	96	6.5				70 - 130	30
4-Methyl-2-pentanone	ND	5.0	76	90	16.9				70 - 130	30
Acetone	ND	5.0	63	73	14.7				70 - 130	30
Acrolein	ND	5.0	93	107	14.0				70 - 130	30
Acrylonitrile	ND	5.0	89	104	15.5				70 - 130	30
Benzene	ND	0.70	89	95	6.5				70 - 130	30
Bromobenzene	ND	1.0	92	98	6.3				70 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank		LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
	Blank	BLK RL								
Bromochloromethane	ND	1.0	88	100	12.8				70 - 130	30
Bromodichloromethane	ND	0.50	89	100	11.6				70 - 130	30
Bromoform	ND	1.0	94	107	12.9				70 - 130	30
Bromomethane	ND	1.0	78	84	7.4				70 - 130	30
Carbon Disulfide	ND	1.0	91	99	8.4				70 - 130	30
Carbon tetrachloride	ND	1.0	89	95	6.5				70 - 130	30
Chlorobenzene	ND	1.0	92	98	6.3				70 - 130	30
Chloroethane	ND	1.0	96	105	9.0				70 - 130	30
Chloroform	ND	1.0	86	93	7.8				70 - 130	30
Chloromethane	ND	1.0	76	85	11.2				70 - 130	30
cis-1,2-Dichloroethene	ND	1.0	90	97	7.5				70 - 130	30
cis-1,3-Dichloropropene	ND	0.40	92	101	9.3				70 - 130	30
Dibromochloromethane	ND	0.50	96	111	14.5				70 - 130	30
Dibromomethane	ND	1.0	89	102	13.6				70 - 130	30
Dichlorodifluoromethane	ND	1.0	98	107	8.8				70 - 130	30
Ethylbenzene	ND	1.0	93	98	5.2				70 - 130	30
Hexachlorobutadiene	ND	0.40	97	100	3.0				70 - 130	30
Isopropylbenzene	ND	1.0	91	95	4.3				70 - 130	30
m&p-Xylene	ND	1.0	91	97	6.4				70 - 130	30
Methyl ethyl ketone	ND	5.0	71	87	20.3				70 - 130	30
Methyl t-butyl ether (MTBE)	ND	1.0	91	107	16.2				70 - 130	30
Methylene chloride	ND	1.0	87	97	10.9				70 - 130	30
Naphthalene	ND	1.0	95	112	16.4				70 - 130	30
n-Butylbenzene	ND	1.0	94	97	3.1				70 - 130	30
n-Propylbenzene	ND	1.0	92	96	4.3				70 - 130	30
o-Xylene	ND	1.0	92	99	7.3				70 - 130	30
p-Isopropyltoluene	ND	1.0	94	96	2.1				70 - 130	30
sec-Butylbenzene	ND	1.0	96	99	3.1				70 - 130	30
Styrene	ND	1.0	92	99	7.3				70 - 130	30
tert-butyl alcohol	ND	10	104	110	5.6				70 - 130	30
tert-Butylbenzene	ND	1.0	91	95	4.3				70 - 130	30
Tetrachloroethene	ND	1.0	90	98	8.5				70 - 130	30
Tetrahydrofuran (THF)	ND	2.5	84	99	16.4				70 - 130	30
Toluene	ND	1.0	90	97	7.5				70 - 130	30
trans-1,2-Dichloroethene	ND	1.0	90	98	8.5				70 - 130	30
trans-1,3-Dichloropropene	ND	0.40	89	99	10.6				70 - 130	30
trans-1,4-dichloro-2-butene	ND	5.0	96	107	10.8				70 - 130	30
Trichloroethene	ND	1.0	92	97	5.3				70 - 130	30
Trichlorofluoromethane	ND	1.0	94	101	7.2				70 - 130	30
Trichlorotrifluoroethane	ND	1.0	93	101	8.2				70 - 130	30
Vinyl chloride	ND	1.0	94	101	7.2				70 - 130	30
% 1,2-dichlorobenzene-d4	100	%	99	101	2.0				70 - 130	30
% Bromofluorobenzene	95	%	100	102	2.0				70 - 130	30
% Dibromofluoromethane	101	%	100	103	3.0				70 - 130	30
% Toluene-d8	99	%	100	99	1.0				70 - 130	30

Comment:

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

Additional 8260 criteria: 10% of LCS/LCSD compounds can be outside of acceptance criteria as long as recovery is 40-160%.

QA/QC Batch 460164 (ug/kg), QC Sample No: CC13978 (CC13976 (50X) , CC13978 (50X))

Volatiles - Soil

1,1,2,2-Tetrachloroethane	ND	3.0	116	111	4.4	107	106	0.9	70 - 130	30
1,2,3-Trichlorobenzene	ND	5.0	109	106	2.8	103	104	1.0	70 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
1,2,3-Trichloropropane	ND	5.0	117	115	1.7	111	110	0.9	70 - 130	30
1,2,4-Trichlorobenzene	ND	5.0	112	105	6.5	110	112	1.8	70 - 130	30
1,2,4-Trimethylbenzene	ND	1.0	108	108	0.0	109	110	0.9	70 - 130	30
1,2-Dibromo-3-chloropropane	ND	5.0	111	108	2.7	100	98	2.0	70 - 130	30
1,2-Dichlorobenzene	ND	5.0	111	110	0.9	109	109	0.0	70 - 130	30
1,3,5-Trimethylbenzene	ND	1.0	108	108	0.0	110	110	0.0	70 - 130	30
1,3-Dichlorobenzene	ND	5.0	112	109	2.7	111	111	0.0	70 - 130	30
1,4-Dichlorobenzene	ND	5.0	112	108	3.6	111	110	0.9	70 - 130	30
2-Chlorotoluene	ND	5.0	109	111	1.8	110	110	0.0	70 - 130	30
2-Isopropyltoluene	ND	5.0	113	112	0.9	114	114	0.0	70 - 130	30
4-Chlorotoluene	ND	5.0	109	109	0.0	110	109	0.9	70 - 130	30
Bromobenzene	ND	5.0	109	109	0.0	107	108	0.9	70 - 130	30
Hexachlorobutadiene	ND	5.0	111	107	3.7	117	118	0.9	70 - 130	30
Isopropylbenzene	ND	1.0	109	110	0.9	109	109	0.0	70 - 130	30
Naphthalene	ND	5.0	108	105	2.8	99	99	0.0	70 - 130	30
n-Butylbenzene	ND	1.0	113	108	4.5	116	116	0.0	70 - 130	30
n-Propylbenzene	ND	1.0	110	110	0.0	111	112	0.9	70 - 130	30
p-Isopropyltoluene	ND	1.0	111	109	1.8	114	114	0.0	70 - 130	30
sec-Butylbenzene	ND	1.0	116	114	1.7	117	117	0.0	70 - 130	30
tert-Butylbenzene	ND	1.0	109	109	0.0	109	110	0.9	70 - 130	30
trans-1,4-dichloro-2-butene	ND	5.0	118	117	0.9	111	111	0.0	70 - 130	30
% 1,2-dichlorobenzene-d4	100	%	99	99	0.0	99	100	1.0	70 - 130	30
% Bromofluorobenzene	97	%	99	99	0.0	99	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%.

QA/QC Batch 460156 (ug/kg), QC Sample No: CC14920 (CC13976, CC13977, CC13978, CC13979, CC13980)

Volatiles - Soil

1,1,1,2-Tetrachloroethane	ND	5.0	102	102	0.0	112	110	1.8	70 - 130	30
1,1,1-Trichloroethane	ND	5.0	105	107	1.9	116	114	1.7	70 - 130	30
1,1,2,2-Tetrachloroethane	ND	3.0	105	104	1.0	109	109	0.0	70 - 130	30
1,1,2-Trichloroethane	ND	5.0	99	97	2.0	108	108	0.0	70 - 130	30
1,1-Dichloroethane	ND	5.0	104	104	0.0	113	112	0.9	70 - 130	30
1,1-Dichloroethene	ND	5.0	105	106	0.9	113	113	0.0	70 - 130	30
1,1-Dichloropropene	ND	5.0	107	108	0.9	119	118	0.8	70 - 130	30
1,2,3-Trichlorobenzene	ND	5.0	97	96	1.0	98	107	8.8	70 - 130	30
1,2,3-Trichloropropane	ND	5.0	103	103	0.0	112	112	0.0	70 - 130	30
1,2,4-Trichlorobenzene	ND	5.0	99	96	3.1	108	113	4.5	70 - 130	30
1,2,4-Trimethylbenzene	ND	1.0	102	102	0.0	111	111	0.0	70 - 130	30
1,2-Dibromo-3-chloropropane	ND	5.0	101	103	2.0	98	105	6.9	70 - 130	30
1,2-Dibromoethane	ND	5.0	100	101	1.0	110	109	0.9	70 - 130	30
1,2-Dichlorobenzene	ND	5.0	102	102	0.0	112	113	0.9	70 - 130	30
1,2-Dichloroethane	ND	5.0	105	104	1.0	117	116	0.9	70 - 130	30
1,2-Dichloropropane	ND	5.0	100	99	1.0	108	108	0.0	70 - 130	30
1,3,5-Trimethylbenzene	ND	1.0	102	102	0.0	111	111	0.0	70 - 130	30
1,3-Dichlorobenzene	ND	5.0	103	101	2.0	114	114	0.0	70 - 130	30
1,3-Dichloropropane	ND	5.0	102	102	0.0	111	108	2.7	70 - 130	30
1,4-Dichlorobenzene	ND	5.0	102	101	1.0	113	114	0.9	70 - 130	30
1,4-dioxane	ND	100	103	103	0.0	107	109	1.9	70 - 130	30
2,2-Dichloropropane	ND	5.0	110	110	0.0	122	121	0.8	70 - 130	30
2-Chlorotoluene	ND	5.0	104	104	0.0	111	112	0.9	70 - 130	30
2-Hexanone	ND	25	76	78	2.6	75	77	2.6	70 - 130	30
2-Isopropyltoluene	ND	5.0	106	106	0.0	107	107	0.0	70 - 130	30

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
4-Chlorotoluene	ND	5.0	102	100	2.0	110	110	0.0	70 - 130	30
4-Methyl-2-pentanone	ND	25	82	84	2.4	85	86	1.2	70 - 130	30
Acetone	ND	10	61	62	1.6	45	45	0.0	70 - 130	30
Acrolein	ND	25	95	97	2.1	83	84	1.2	70 - 130	30
Acrylonitrile	ND	5.0	97	100	3.0	99	98	1.0	70 - 130	30
Benzene	ND	1.0	101	102	1.0	110	110	0.0	70 - 130	30
Bromobenzene	ND	5.0	102	101	1.0	108	109	0.9	70 - 130	30
Bromochloromethane	ND	5.0	105	105	0.0	115	114	0.9	70 - 130	30
Bromodichloromethane	ND	5.0	101	102	1.0	112	111	0.9	70 - 130	30
Bromoform	ND	5.0	98	97	1.0	105	103	1.9	70 - 130	30
Bromomethane	ND	5.0	122	120	1.7	85	97	13.2	70 - 130	30
Carbon Disulfide	ND	5.0	106	106	0.0	107	106	0.9	70 - 130	30
Carbon tetrachloride	ND	5.0	92	107	15.1	101	113	11.2	70 - 130	30
Chlorobenzene	ND	5.0	104	104	0.0	113	112	0.9	70 - 130	30
Chloroethane	ND	5.0	116	114	1.7	124	123	0.8	70 - 130	30
Chloroform	ND	5.0	104	103	1.0	113	111	1.8	70 - 130	30
Chloromethane	ND	5.0	94	95	1.1	100	99	1.0	70 - 130	30
cis-1,2-Dichloroethene	ND	5.0	106	105	0.9	113	109	3.6	70 - 130	30
cis-1,3-Dichloropropene	ND	5.0	104	103	1.0	114	112	1.8	70 - 130	30
Dibromochloromethane	ND	3.0	104	105	1.0	111	112	0.9	70 - 130	30
Dibromomethane	ND	5.0	99	100	1.0	111	110	0.9	70 - 130	30
Dichlorodifluoromethane	ND	5.0	121	123	1.6	134	132	1.5	70 - 130	30
Ethylbenzene	ND	1.0	104	104	0.0	114	113	0.9	70 - 130	30
Hexachlorobutadiene	ND	5.0	102	104	1.9	117	117	0.0	70 - 130	30
Isopropylbenzene	ND	1.0	104	105	1.0	111	110	0.9	70 - 130	30
m&p-Xylene	ND	2.0	103	103	0.0	114	113	0.9	70 - 130	30
Methyl ethyl ketone	ND	5.0	79	79	0.0	82	79	3.7	70 - 130	30
Methyl t-butyl ether (MTBE)	ND	1.0	103	102	1.0	102	101	1.0	70 - 130	30
Methylene chloride	ND	5.0	88	87	1.1	97	94	3.1	70 - 130	30
Naphthalene	ND	5.0	97	97	0.0	91	102	11.4	70 - 130	30
n-Butylbenzene	ND	1.0	104	105	1.0	116	117	0.9	70 - 130	30
n-Propylbenzene	ND	1.0	105	106	0.9	112	112	0.0	70 - 130	30
o-Xylene	ND	2.0	103	103	0.0	114	112	1.8	70 - 130	30
p-Isopropyltoluene	ND	1.0	105	105	0.0	114	114	0.0	70 - 130	30
sec-Butylbenzene	ND	1.0	109	110	0.9	118	117	0.9	70 - 130	30
Styrene	ND	5.0	102	101	1.0	113	112	0.9	70 - 130	30
tert-butyl alcohol	ND	100	105	101	3.9	91	101	10.4	70 - 130	30
tert-Butylbenzene	ND	1.0	104	104	0.0	111	110	0.9	70 - 130	30
Tetrachloroethene	ND	5.0	103	104	1.0	116	117	0.9	70 - 130	30
Tetrahydrofuran (THF)	ND	5.0	98	97	1.0	96	95	1.0	70 - 130	30
Toluene	ND	1.0	101	101	0.0	112	111	0.9	70 - 130	30
trans-1,2-Dichloroethene	ND	5.0	105	106	0.9	116	112	3.5	70 - 130	30
trans-1,3-Dichloropropene	ND	5.0	101	100	1.0	113	111	1.8	70 - 130	30
trans-1,4-dichloro-2-butene	ND	5.0	109	108	0.9	103	104	1.0	70 - 130	30
Trichloroethene	ND	5.0	106	108	1.9	117	117	0.0	70 - 130	30
Trichlorofluoromethane	ND	5.0	115	115	0.0	114	112	1.8	70 - 130	30
Trichlorotrifluoroethane	ND	5.0	105	107	1.9	107	105	1.9	70 - 130	30
Vinyl chloride	ND	5.0	111	110	0.9	115	113	1.8	70 - 130	30
% 1,2-dichlorobenzene-d4	100	%	101	99	2.0	100	101	1.0	70 - 130	30
% Bromofluorobenzene	97	%	100	98	2.0	101	100	1.0	70 - 130	30
% Dibromofluoromethane	97	%	96	95	1.0	97	97	0.0	70 - 130	30
% Toluene-d8	99	%	100	99	1.0	100	100	0.0	70 - 130	30

l,m

m

QA/QC Data

SDG I.D.: GCC13976

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
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Comment:

Additional 8260 criteria: 10% of LCS/LCSD compounds can be outside of acceptance criteria as long as recovery is 40-160%.

l = 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

December 19, 2018

Wednesday, December 19, 2018

Criteria: NY: 375, 375GWP, 375RRS, 375RS, GW

State: NY

Sample Criteria Exceedances Report

GCC13976 - EBC

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	Criteria	RL	Analysis Units
CC13976	\$8270SMRDP	Benz(a)anthracene	NY / 375-6.8 Semivolatiles / Ground Water Protection	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(b)fluoranthene	NY / 375-6.8 Semivolatiles / Ground Water Protection	3000	260	1700	1700	1700	ug/Kg
CC13976	\$8270SMRDP	Chrysene	NY / 375-6.8 Semivolatiles / Ground Water Protection	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(k)fluoranthene	NY / 375-6.8 Semivolatiles / Ground Water Protection	2700	260	1700	1700	1700	ug/Kg
CC13976	\$8270SMRDP	Dibenz(a,h)anthracene	NY / 375-6.8 Semivolatiles / Residential	720	190	330	330	330	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)anthracene	NY / 375-6.8 Semivolatiles / Residential	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)pyrene	NY / 375-6.8 Semivolatiles / Residential	3200	190	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Indeno(1,2,3-cd)pyrene	NY / 375-6.8 Semivolatiles / Residential	2100	260	500	500	500	ug/Kg
CC13976	\$8270SMRDP	Benzo(b)fluoranthene	NY / 375-6.8 Semivolatiles / Residential	3000	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Chrysene	NY / 375-6.8 Semivolatiles / Residential	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(k)fluoranthene	NY / 375-6.8 Semivolatiles / Residential	2700	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Indeno(1,2,3-cd)pyrene	NY / 375-6.8 Semivolatiles / Residential Restricted	2100	260	500	500	500	ug/Kg
CC13976	\$8270SMRDP	Dibenz(a,h)anthracene	NY / 375-6.8 Semivolatiles / Residential Restricted	720	190	330	330	330	ug/Kg
CC13976	\$8270SMRDP	Benzo(b)fluoranthene	NY / 375-6.8 Semivolatiles / Residential Restricted	3000	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)anthracene	NY / 375-6.8 Semivolatiles / Residential Restricted	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)pyrene	NY / 375-6.8 Semivolatiles / Residential Restricted	3200	190	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Indeno(1,2,3-cd)pyrene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	2100	260	500	500	500	ug/Kg
CC13976	\$8270SMRDP	Benzo(b)fluoranthene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	3000	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)pyrene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	3200	190	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Chrysene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Dibenz(a,h)anthracene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	720	190	330	330	330	ug/Kg
CC13976	\$8270SMRDP	Benzo(a)anthracene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	3400	260	1000	1000	1000	ug/Kg
CC13976	\$8270SMRDP	Benzo(k)fluoranthene	NY / 375-6.8 Semivolatiles / Unrestricted Use Soil	2700	260	800	800	800	ug/Kg
CC13976	AS-SM	Arsenic	NY / 375-6.8 Metals / Unrestricted Use Soil	15.8	0.75	13	13	13	mg/Kg
CC13976	CD-SM	Cadmium	NY / 375-6.8 Metals / Residential	3.44	0.37	2.5	2.5	2.5	mg/Kg
CC13976	CD-SM	Cadmium	NY / 375-6.8 Metals / Unrestricted Use Soil	3.44	0.37	2.5	2.5	2.5	mg/Kg
CC13976	CR-SM	Chromium	NY / 375-6.8 Metals / Unrestricted Use Soil	46.3	0.37	30			mg/Kg
CC13976	CU-SM	Copper	NY / 375-6.8 Metals / Residential	308	7.5	270	270	270	mg/kg
CC13976	CU-SM	Copper	NY / 375-6.8 Metals / Residential Restricted	308	7.5	270	270	270	mg/kg
CC13976	CU-SM	Copper	NY / 375-6.8 Metals / Unrestricted Use Soil	308	7.5	50	50	50	mg/kg
CC13976	HG-SM	Mercury	NY / 375-6.8 Metals / Ground Water Protection	1.29	0.15	0.73	0.73	0.73	mg/Kg
CC13976	HG-SM	Mercury	NY / 375-6.8 Metals / Residential	1.29	0.15	0.81	0.81	0.81	mg/Kg
CC13976	HG-SM	Mercury	NY / 375-6.8 Metals / Residential Restricted	1.29	0.15	0.81	0.81	0.81	mg/Kg
CC13976	HG-SM	Mercury	NY / 375-6.8 Metals / Unrestricted Use Soil	1.29	0.15	0.18	0.18	0.18	mg/Kg
CC13976	NI-SM	Nickel	NY / 375-6.8 Metals / Unrestricted Use Soil	38.3	0.37	30	30	30	mg/Kg
CC13976	PB-SMDP	Lead	NY / 375-6.8 Metals / Ground Water Protection	662	7.5	450	450	450	mg/Kg
CC13976	PB-SMDP	Lead	NY / 375-6.8 Metals / Residential	662	7.5	400	400	400	mg/Kg
CC13976	PB-SMDP	Lead	NY / 375-6.8 Metals / Residential Restricted	662	7.5	400	400	400	mg/Kg
CC13976	PB-SMDP	Lead	NY / 375-6.8 Metals / Unrestricted Use Soil	662	7.5	63	63	63	mg/Kg
CC13976	ZN-SMDP	Zinc	NY / 375-6.8 Metals / Unrestricted Use Soil	720	7.5	109	109	109	mg/Kg
CC13977	CD-SM	Cadmium	NY / 375-6.8 Metals / Residential	2.61	0.40	2.5	2.5	2.5	mg/Kg

Wednesday, December 19, 2018

Criteria: NY: 375, 375GWP, 375RRS, 375RS, GW

State: NY

Sample Criteria Exceedances Report

GCC13976 - EBC

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	RL Criteria	Analysis Units
CC13977	CD-SM	Cadmium	NY / 375-6.8 Metals / Unrestricted Use Soil	2.61	0.40	2.5	2.5	mg/Kg
CC13977	CU-SM	Copper	NY / 375-6.8 Metals / Unrestricted Use Soil	159	0.8	50	50	mg/kg
CC13977	HG-SM	Mercury	NY / 375-6.8 Metals / Ground Water Protection	13.0	1.4	0.73	0.73	mg/Kg
CC13977	HG-SM	Mercury	NY / 375-6.8 Metals / Residential	13.0	1.4	0.81	0.81	mg/Kg
CC13977	HG-SM	Mercury	NY / 375-6.8 Metals / Residential Restricted	13.0	1.4	0.81	0.81	mg/Kg
CC13977	HG-SM	Mercury	NY / 375-6.8 Metals / Unrestricted Use Soil	13.0	1.4	0.18	0.18	mg/Kg
CC13977	PB-SMDP	Lead	NY / 375-6.8 Metals / Ground Water Protection	959	8.0	450	450	mg/Kg
CC13977	PB-SMDP	Lead	NY / 375-6.8 Metals / Residential	959	8.0	400	400	mg/Kg
CC13977	PB-SMDP	Lead	NY / 375-6.8 Metals / Residential Restricted	959	8.0	400	400	mg/Kg
CC13977	PB-SMDP	Lead	NY / 375-6.8 Metals / Unrestricted Use Soil	959	8.0	63	63	mg/Kg
CC13977	ZN-SMDP	Zinc	NY / 375-6.8 Metals / Unrestricted Use Soil	877	8.0	109	109	mg/Kg
CC13978	CR-SM	Chromium	NY / 375-6.8 Metals / Unrestricted Use Soil	34.8	0.37	30		mg/Kg
CC13978	CU-SM	Copper	NY / 375-6.8 Metals / Unrestricted Use Soil	63.9	0.7	50	50	mg/kg
CC13978	HG-SM	Mercury	NY / 375-6.8 Metals / Unrestricted Use Soil	0.53	0.14	0.18	0.18	mg/Kg
CC13978	NI-SM	Nickel	NY / 375-6.8 Metals / Unrestricted Use Soil	43.2	0.37	30	30	mg/Kg
CC13978	PB-SMDP	Lead	NY / 375-6.8 Metals / Unrestricted Use Soil	382	7.4	63	63	mg/Kg
CC13978	ZN-SMDP	Zinc	NY / 375-6.8 Metals / Unrestricted Use Soil	195	7.4	109	109	mg/Kg
CC13979	HG-SM	Mercury	NY / 375-6.8 Metals / Unrestricted Use Soil	0.24	0.13	0.18	0.18	mg/Kg
CC13981	\$8260DP25R	1,2-Dibromoethane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.0006	0.0006	ug/L
CC13981	\$8260DP25R	1,2,3-Trichloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.04	0.04	ug/L
CC13981	\$8260DP25R	1,2-Dibromo-3-chloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.50	0.04	0.04	ug/L
CC13982	\$8260DP25R	1,2,3-Trichloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.04	0.04	ug/L
CC13982	\$8260DP25R	1,2-Dibromo-3-chloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.50	0.04	0.04	ug/L
CC13982	\$8260DP25R	1,2-Dibromoethane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.0006	0.0006	ug/L
CC13983	\$8260DP25R	1,2-Dibromoethane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.0006	0.0006	ug/L
CC13983	\$8260DP25R	1,2,3-Trichloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.25	0.04	0.04	ug/L
CC13983	\$8260DP25R	1,2-Dibromo-3-chloropropane	NY / TOGS - Water Quality / GA Criteria	ND	0.50	0.04	0.04	ug/L

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 exceedances. All efforts are made to ensure the accuracy of the data (obtained from appropriate agencies). A lack of exceedance 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
Tel. (860) 645-1102 Fax (860) 645-0823



Analysis Comments

December 19, 2018

SDG I.D.: GCC13976

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

SVOA Narration

CHEM19 12/13/18-3: CC13976, CC13977, CC13978, CC13979, CC13980

The following Initial Calibration compounds did not meet RSD% criteria: 2,4-Dinitrophenol 26% (20%), 4,6-Dinitro-2-methylphenol 22% (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-Nitrophenol 0.054 (0.1), Hexachlorobenzene 0.090 (0.1)

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

The following Continuing Calibration compounds did not meet recommended response factors: 2-Nitrophenol 0.056 (0.1), Hexachlorobenzene 0.086 (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

CHEM02 12/14/18-2: CC13982

The following Initial Calibration compounds did not meet RSD% criteria: Acetone 29% (20%)

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

The following Initial Calibration compounds did not meet recommended response factors: 1,2-Dibromo-3-chloropropane 0.028 (0.05), 2-Hexanone 0.085 (0.1), Acetone 0.057 (0.1), Acrolein 0.021 (0.05), Bromoform 0.084 (0.1), Methyl ethyl ketone 0.070 (0.1), Tetrahydrofuran (THF) 0.042 (0.05)

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

The following Continuing Calibration compounds did not meet % deviation criteria: Bromomethane 35%L (30%)

The following Continuing Calibration compounds did not meet Maximum % deviation criteria: None.

The following Continuing Calibration compounds did not meet recommended response factors: 1,1,2,2-Tetrachloroethane 0.242 (0.3), 1,2-Dibromo-3-chloropropane 0.029 (0.05), Acrolein 0.019 (0.05), Acrylonitrile 0.049 (0.05), Bromoform 0.083 (0.1), Tetrahydrofuran (THF) 0.038 (0.05)

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%.

CHEM02 12/15/18-1: CC13981, CC13983

The following Initial Calibration compounds did not meet RSD% criteria: Acetone 29% (20%)

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

The following Initial Calibration compounds did not meet recommended response factors: 1,2-Dibromo-3-chloropropane 0.028 (0.05), 2-Hexanone 0.085 (0.1), Acetone 0.057 (0.1), Acrolein 0.021 (0.05), Bromoform 0.084 (0.1), Methyl ethyl ketone 0.070 (0.1), Tetrahydrofuran (THF) 0.042 (0.05)

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

The following Continuing Calibration compounds did not meet % deviation criteria: Bromomethane 33%L (30%)

The following Continuing Calibration compounds did not meet Maximum % deviation criteria: None.

The following Continuing Calibration compounds did not meet recommended response factors: 1,1,2,2-Tetrachloroethane 0.242 (0.3), 1,2-Dibromo-3-chloropropane 0.030 (0.05), Acrolein 0.021 (0.05), Bromoform 0.089 (0.1), Tetrahydrofuran (THF) 0.037 (0.05)

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%.



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

December 19, 2018

SDG I.D.: GCC13976

CHEM03 12/14/18-2: CC13976, CC13977, CC13978, CC13979, CC13980

The following Initial Calibration compounds did not meet RSD% criteria: 1,2,3-Trichlorobenzene 25% (20%), Acetone 36% (20%), Methyl Ethyl Ketone 21% (20%), Naphthalene 33% (20%)

The following Initial Calibration compounds did not meet maximum RSD% criteria: 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%.

CHEM03 12/15/18-1: CC13976, CC13978

The following Initial Calibration compounds did not meet RSD% criteria: 1,2,3-Trichlorobenzene 25% (20%), Naphthalene 33% (20%)

The following Initial Calibration compounds did not meet maximum RSD% criteria: 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

December 19, 2018

SDG I.D.: GCC13976

The samples in this delivery group were received at 3.6°C.
(Note acceptance criteria for relevant matrices is above freezing up to 6°C)



NY/NJ CHAIN OF CUSTODY RECORD

587 East Middle Turnpike, P.O. Box 370, Manchester, CT 06040
 Email: info@phoenixlabs.com Fax (860) 645-0823
Client Services (860) 645-8726

Customer: Environmental Business Consultants
Address: 1808 Middle Country Road
 Ridge, NY 11961

Project: 188 East 135th Street, Grand, NJ
Report to: Environmental Business Consultants
Invoice to: Environmental Business Consultants

Project P.O.:

This section MUST be completed with Bottle Quantities.

Coolant: Yes No
 IPK ICE No
 Temp: 31°C Pg 1 of 1

Contact Options:
 Fax:
 Phone: 631-504-6000
 Email: F.I.C.

Sampler's Signature	Client Sample - Information - Identification	Analysis Request
David Rukk	Date: 12-12-18	VOCs 8469 SVCs 8470 TAL Metals

PHOENIX USE ONLY	Customer Sample Identification	Sample Matrix	Date Sampled	Time Sampled
139710	S61 (10-12')	S	12-12-18	8:15
139717	S62 (7-9')	S		11:45
139718	S63 (7-9')	S		10:30
139719	S64 (8-10')	S		9:45
139780	S65 (8-10')	S		9:00
139781	GW1	GW		8:40
139782	GW2	GW		10:00
139783	GW3	GW		10:45

Relinquished by:	Accepted by:	Date:	Time:
David Rukk	[Signature]	12-13-18	13:00
[Signature]	[Signature]	12-13-18	14:30

Turnaround:
 1 Day*
 2 Days*
 3 Days*
 5 Days
 10 Days
 Other
 *SURCHARGE APPLIES

NJ
 Res. Criteria
 Non-Res. Criteria
 Impact to GW Soil Cleanup Criteria
 GW Criteria

NY
 NY 375 GWP
 NY375 Unrestricted Use Soil
 NY375 Residential Soil
 Restricted/Residential Commercial
 Industrial

Data Format
 Phoenix Std Report
 Excel
 PDF
 GIS/Key
 EQUIS
 NJ Hazsite EDD
 NY EZ EDD (ASP)
 Other

Data Package
 NJ Reduced Deliv.*
 NY Enhanced (ASP B)*
 Other

State where samples were collected:
 NY

Comments, Special Requirements or Regulations:



Friday, December 14, 2018

Attn: Mr. Charles B. Sosik, P.G.
Environmental Business Consultants
1808 Middle Country Rd
Ridge NY 11961-2406

Project ID: 188 EAST 135TH ST BRONX
Sample ID#s: CC13973 - CC13975

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.

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

If you have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext. 200.

Sincerely yours,

A handwritten signature in black ink that reads "Phyllis Shiller". The signature is written in a cursive style.

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
UT Lab Registration #CT00007
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

Analysis Report
 December 14, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: AIR
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:
 Canister Id: 23346

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date Time
 12/12/18 11:56
 12/13/18 16:30

Laboratory Data

SDG ID: GCC13973
 Phoenix ID: CC13973

Project ID: 188 EAST 135TH ST BRONX
 Client ID: SV 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution	
<u>Volatiles (TO15)</u>								
1,1,1,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1	1
1,1,1-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1	
1,1,2,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1	
1,1,2-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1	
1,1-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1	
1,1-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1	
1,2,4-Trichlorobenzene	ND	0.135	ND	1.00	12/13/18	KCA	1	
1,2,4-Trimethylbenzene	0.591	0.204	2.90	1.00	12/13/18	KCA	1	
1,2-Dibromoethane(EDB)	ND	0.130	ND	1.00	12/13/18	KCA	1	
1,2-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1	
1,2-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1	
1,2-dichloropropane	ND	0.217	ND	1.00	12/13/18	KCA	1	
1,2-Dichlorotetrafluoroethane	ND	0.143	ND	1.00	12/13/18	KCA	1	
1,3,5-Trimethylbenzene	0.288	0.204	1.41	1.00	12/13/18	KCA	1	
1,3-Butadiene	ND	0.452	ND	1.00	12/13/18	KCA	1	
1,3-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1	
1,4-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1	
1,4-Dioxane	ND	0.278	ND	1.00	12/13/18	KCA	1	
2-Hexanone(MBK)	ND	0.244	ND	1.00	12/13/18	KCA	1	1
4-Ethyltoluene	0.877	0.204	4.31	1.00	12/13/18	KCA	1	1
4-Isopropyltoluene	ND	0.182	ND	1.00	12/13/18	KCA	1	1
4-Methyl-2-pentanone(MIBK)	ND	0.244	ND	1.00	12/13/18	KCA	1	
Acetone	121	2.11	287	5.01	12/14/18	KCA	5	
Acrylonitrile	ND	0.461	ND	1.00	12/13/18	KCA	1	
Benzene	1.67	0.313	5.33	1.00	12/13/18	KCA	1	
Benzyl chloride	ND	0.193	ND	1.00	12/13/18	KCA	1	

Client ID: SV 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
Bromodichloromethane	ND	0.149	ND	1.00	12/13/18	KCA	1
Bromoform	ND	0.097	ND	1.00	12/13/18	KCA	1
Bromomethane	ND	0.258	ND	1.00	12/13/18	KCA	1
Carbon Disulfide	4.73	0.321	14.7	1.00	12/13/18	KCA	1
Carbon Tetrachloride	ND	0.032	ND	0.20	12/13/18	KCA	1
Chlorobenzene	ND	0.217	ND	1.00	12/13/18	KCA	1
Chloroethane	ND	0.379	ND	1.00	12/13/18	KCA	1
Chloroform	ND	0.205	ND	1.00	12/13/18	KCA	1
Chloromethane	ND	0.485	ND	1.00	12/13/18	KCA	1
Cis-1,2-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1
cis-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Cyclohexane	2.49	0.291	8.57	1.00	12/13/18	KCA	1
Dibromochloromethane	ND	0.118	ND	1.00	12/13/18	KCA	1
Dichlorodifluoromethane	0.295	0.202	1.46	1.00	12/13/18	KCA	1
Ethanol	14.6	0.531	27.5	1.00	12/13/18	KCA	1
Ethyl acetate	ND	0.278	ND	1.00	12/13/18	KCA	1
Ethylbenzene	0.926	0.230	4.02	1.00	12/13/18	KCA	1
Heptane	10.2	0.244	41.8	1.00	12/13/18	KCA	1
Hexachlorobutadiene	ND	0.094	ND	1.00	12/13/18	KCA	1
Hexane	22.0	0.284	77.5	1.00	12/13/18	KCA	1
Isopropylalcohol	ND	0.407	ND	1.00	12/13/18	KCA	1
Isopropylbenzene	ND	0.204	ND	1.00	12/13/18	KCA	1
m,p-Xylene	3.47	0.230	15.1	1.00	12/13/18	KCA	1
Methyl Ethyl Ketone	7.46	0.339	22.0	1.00	12/13/18	KCA	1
Methyl tert-butyl ether(MTBE)	ND	0.278	ND	1.00	12/13/18	KCA	1
Methylene Chloride	18.7	0.864	64.9	3.00	12/13/18	KCA	1
n-Butylbenzene	ND	0.182	ND	1.00	12/13/18	KCA	1
o-Xylene	1.23	0.230	5.34	1.00	12/13/18	KCA	1
Propylene	ND	0.581	ND	1.00	12/13/18	KCA	1
sec-Butylbenzene	ND	0.182	ND	1.00	12/13/18	KCA	1
Styrene	0.359	0.235	1.53	1.00	12/13/18	KCA	1
Tetrachloroethene	0.553	0.037	3.75	0.25	12/13/18	KCA	1
Tetrahydrofuran	ND	0.339	ND	1.00	12/13/18	KCA	1
Toluene	4.64	0.266	17.5	1.00	12/13/18	KCA	1
Trans-1,2-Dichloroethene	ND	0.252	ND	1.00	12/13/18	KCA	1
trans-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Trichloroethene	ND	0.037	ND	0.20	12/13/18	KCA	1
Trichlorofluoromethane	ND	0.178	ND	1.00	12/13/18	KCA	1
Trichlorotrifluoroethane	ND	0.131	ND	1.00	12/13/18	KCA	1
Vinyl Chloride	ND	0.078	ND	0.20	12/13/18	KCA	1
QA/QC Surrogates							
% Bromofluorobenzene	88	%	88	%	12/13/18	KCA	1

Client ID: SV 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
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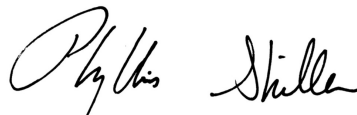
1 = This parameter is not certified by the primary accrediting authority (NY NELAC) for this matrix. NY NELAC does not offer certification for all parameters at this time.

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 (preceeded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

Comments:

If there are any questions regarding this data, please call Phoenix Client Services.
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Phyllis Shiller, Laboratory Director

December 14, 2018

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report
 December 14, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: AIR
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:
 Canister Id: 11290

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date Time
 12/12/18 12:32
 12/13/18 16:30

Laboratory Data

SDG ID: GCC13973
 Phoenix ID: CC13974

Project ID: 188 EAST 135TH ST BRONX
 Client ID: SV 2

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
<u>Volatiles (TO15)</u>							
1,1,1,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1
1,1,1-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1
1,1,2,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1
1,1,2-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1
1,1-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1
1,1-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1
1,2,4-Trichlorobenzene	ND	0.135	ND	1.00	12/13/18	KCA	1
1,2,4-Trimethylbenzene	0.512	0.204	2.52	1.00	12/13/18	KCA	1
1,2-Dibromoethane(EDB)	ND	0.130	ND	1.00	12/13/18	KCA	1
1,2-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,2-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1
1,2-dichloropropane	ND	0.217	ND	1.00	12/13/18	KCA	1
1,2-Dichlorotetrafluoroethane	ND	0.143	ND	1.00	12/13/18	KCA	1
1,3,5-Trimethylbenzene	ND	0.204	ND	1.00	12/13/18	KCA	1
1,3-Butadiene	4.02	0.452	8.89	1.00	12/13/18	KCA	1
1,3-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,4-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,4-Dioxane	ND	0.278	ND	1.00	12/13/18	KCA	1
2-Hexanone(MBK)	ND	0.244	ND	1.00	12/13/18	KCA	1
4-Ethyltoluene	0.764	0.204	3.75	1.00	12/13/18	KCA	1
4-Isopropyltoluene	ND	0.182	ND	1.00	12/13/18	KCA	1
4-Methyl-2-pentanone(MIBK)	0.557	0.244	2.28	1.00	12/13/18	KCA	1
Acetone	104	2.11	247	5.01	12/14/18	KCA	5
Acrylonitrile	ND	0.461	ND	1.00	12/13/18	KCA	1
Benzene	0.869	0.313	2.77	1.00	12/13/18	KCA	1
Benzyl chloride	ND	0.193	ND	1.00	12/13/18	KCA	1

Client ID: SV 2

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
Bromodichloromethane	ND	0.149	ND	1.00	12/13/18	KCA	1
Bromoform	ND	0.097	ND	1.00	12/13/18	KCA	1
Bromomethane	ND	0.258	ND	1.00	12/13/18	KCA	1
Carbon Disulfide	1.65	0.321	5.13	1.00	12/13/18	KCA	1
Carbon Tetrachloride	ND	0.032	ND	0.20	12/13/18	KCA	1
Chlorobenzene	ND	0.217	ND	1.00	12/13/18	KCA	1
Chloroethane	ND	0.379	ND	1.00	12/13/18	KCA	1
Chloroform	ND	0.205	ND	1.00	12/13/18	KCA	1
Chloromethane	ND	0.485	ND	1.00	12/13/18	KCA	1
Cis-1,2-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1
cis-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Cyclohexane	2.94	0.291	10.1	1.00	12/13/18	KCA	1
Dibromochloromethane	ND	0.118	ND	1.00	12/13/18	KCA	1
Dichlorodifluoromethane	0.342	0.202	1.69	1.00	12/13/18	KCA	1
Ethanol	22.3	0.531	42.0	1.00	12/13/18	KCA	1
Ethyl acetate	ND	0.278	ND	1.00	12/13/18	KCA	1
Ethylbenzene	0.577	0.230	2.50	1.00	12/13/18	KCA	1
Heptane	0.936	0.244	3.83	1.00	12/13/18	KCA	1
Hexachlorobutadiene	ND	0.094	ND	1.00	12/13/18	KCA	1
Hexane	1.63	S 0.284	5.74	1.00	12/13/18	KCA	1
Isopropylalcohol	0.710	0.407	1.74	1.00	12/13/18	KCA	1
Isopropylbenzene	ND	0.204	ND	1.00	12/13/18	KCA	1
m,p-Xylene	2.37	0.230	10.3	1.00	12/13/18	KCA	1
Methyl Ethyl Ketone	3.90	0.339	11.5	1.00	12/13/18	KCA	1
Methyl tert-butyl ether(MTBE)	ND	0.278	ND	1.00	12/13/18	KCA	1
Methylene Chloride	3.52	S 0.864	12.2	3.00	12/13/18	KCA	1
n-Butylbenzene	ND	0.182	ND	1.00	12/13/18	KCA	1
o-Xylene	0.779	0.230	3.38	1.00	12/13/18	KCA	1
Propylene	97.4	2.91	168	5.01	12/14/18	KCA	5
sec-Butylbenzene	ND	0.182	ND	1.00	12/13/18	KCA	1
Styrene	0.280	0.235	1.19	1.00	12/13/18	KCA	1
Tetrachloroethene	0.091	0.037	0.62	0.25	12/13/18	KCA	1
Tetrahydrofuran	2.69	0.339	7.93	1.00	12/13/18	KCA	1
Toluene	2.80	0.266	10.5	1.00	12/13/18	KCA	1
Trans-1,2-Dichloroethene	ND	0.252	ND	1.00	12/13/18	KCA	1
trans-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Trichloroethene	ND	0.037	ND	0.20	12/13/18	KCA	1
Trichlorofluoromethane	ND	0.178	ND	1.00	12/13/18	KCA	1
Trichlorotrifluoroethane	ND	0.131	ND	1.00	12/13/18	KCA	1
Vinyl Chloride	0.190	0.078	0.49	0.20	12/13/18	KCA	1
<u>QA/QC Surrogates</u>							
% Bromofluorobenzene	88	%	88	%	12/13/18	KCA	1

Client ID: SV 2

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
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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 (preceeded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

Comments:

S - Laboratory solvent, contamination is possible.

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

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Phyllis Shiller, Laboratory Director

December 14, 2018

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report
 December 14, 2018

FOR: Attn: Mr. Charles B. Sosik, P.G.
 Environmental Business Consultants
 1808 Middle Country Rd
 Ridge NY 11961-2406

Sample Information

Matrix: AIR
 Location Code: EBC
 Rush Request: 72 Hour
 P.O.#:
 Canister Id: 19425

Custody Information

Collected by: DR
 Received by: SW
 Analyzed by: see "By" below

Date Time
 12/12/18 12:25
 12/13/18 16:30

Laboratory Data

SDG ID: GCC13973
 Phoenix ID: CC13975

Project ID: 188 EAST 135TH ST BRONX
 Client ID: SS 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
<u>Volatiles (TO15)</u>							
1,1,1,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1
1,1,1-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1
1,1,2,2-Tetrachloroethane	ND	0.146	ND	1.00	12/13/18	KCA	1
1,1,2-Trichloroethane	ND	0.183	ND	1.00	12/13/18	KCA	1
1,1-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1
1,1-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1
1,2,4-Trichlorobenzene	ND	0.135	ND	1.00	12/13/18	KCA	1
1,2,4-Trimethylbenzene	0.979	0.204	4.81	1.00	12/13/18	KCA	1
1,2-Dibromoethane(EDB)	ND	0.130	ND	1.00	12/13/18	KCA	1
1,2-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,2-Dichloroethane	ND	0.247	ND	1.00	12/13/18	KCA	1
1,2-dichloropropane	ND	0.217	ND	1.00	12/13/18	KCA	1
1,2-Dichlorotetrafluoroethane	ND	0.143	ND	1.00	12/13/18	KCA	1
1,3,5-Trimethylbenzene	0.350	0.204	1.72	1.00	12/13/18	KCA	1
1,3-Butadiene	ND	0.452	ND	1.00	12/13/18	KCA	1
1,3-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,4-Dichlorobenzene	ND	0.166	ND	1.00	12/13/18	KCA	1
1,4-Dioxane	ND	0.278	ND	1.00	12/13/18	KCA	1
2-Hexanone(MBK)	ND	0.244	ND	1.00	12/13/18	KCA	1
4-Ethyltoluene	1.25	0.204	6.14	1.00	12/13/18	KCA	1
4-Isopropyltoluene	ND	0.182	ND	1.00	12/13/18	KCA	1
4-Methyl-2-pentanone(MIBK)	0.352	0.244	1.44	1.00	12/13/18	KCA	1
Acetone	62.5	2.11	148	5.01	12/14/18	KCA	5
Acrylonitrile	ND	0.461	ND	1.00	12/13/18	KCA	1
Benzene	1.84	0.313	5.87	1.00	12/13/18	KCA	1
Benzyl chloride	ND	0.193	ND	1.00	12/13/18	KCA	1

Client ID: SS 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
Bromodichloromethane	ND	0.149	ND	1.00	12/13/18	KCA	1
Bromoform	ND	0.097	ND	1.00	12/13/18	KCA	1
Bromomethane	ND	0.258	ND	1.00	12/13/18	KCA	1
Carbon Disulfide	1.17	0.321	3.64	1.00	12/13/18	KCA	1
Carbon Tetrachloride	0.053	0.032	0.33	0.20	12/13/18	KCA	1
Chlorobenzene	ND	0.217	ND	1.00	12/13/18	KCA	1
Chloroethane	ND	0.379	ND	1.00	12/13/18	KCA	1
Chloroform	ND	0.205	ND	1.00	12/13/18	KCA	1
Chloromethane	ND	0.485	ND	1.00	12/13/18	KCA	1
Cis-1,2-Dichloroethene	ND	0.051	ND	0.20	12/13/18	KCA	1
cis-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Cyclohexane	0.310	0.291	1.07	1.00	12/13/18	KCA	1
Dibromochloromethane	ND	0.118	ND	1.00	12/13/18	KCA	1
Dichlorodifluoromethane	0.304	0.202	1.50	1.00	12/13/18	KCA	1
Ethanol	10.5	0.531	19.8	1.00	12/13/18	KCA	1
Ethyl acetate	ND	0.278	ND	1.00	12/13/18	KCA	1
Ethylbenzene	1.26	0.230	5.47	1.00	12/13/18	KCA	1
Heptane	0.797	0.244	3.26	1.00	12/13/18	KCA	1
Hexachlorobutadiene	ND	0.094	ND	1.00	12/13/18	KCA	1
Hexane	0.375	S 0.284	1.32	1.00	12/13/18	KCA	1
Isopropylalcohol	1.17	0.407	2.87	1.00	12/13/18	KCA	1
Isopropylbenzene	ND	0.204	ND	1.00	12/13/18	KCA	1
m,p-Xylene	4.22	0.230	18.3	1.00	12/13/18	KCA	1
Methyl Ethyl Ketone	4.24	0.339	12.5	1.00	12/13/18	KCA	1
Methyl tert-butyl ether(MTBE)	ND	0.278	ND	1.00	12/13/18	KCA	1
Methylene Chloride	ND	0.864	ND	3.00	12/13/18	KCA	1
n-Butylbenzene	0.223	0.182	1.22	1.00	12/13/18	KCA	1
o-Xylene	1.49	0.230	6.47	1.00	12/13/18	KCA	1
Propylene	1.62	0.581	2.79	1.00	12/13/18	KCA	1
sec-Butylbenzene	ND	0.182	ND	1.00	12/13/18	KCA	1
Styrene	0.676	0.235	2.88	1.00	12/13/18	KCA	1
Tetrachloroethene	0.846	0.037	5.73	0.25	12/13/18	KCA	1
Tetrahydrofuran	ND	0.339	ND	1.00	12/13/18	KCA	1
Toluene	5.74	0.266	21.6	1.00	12/13/18	KCA	1
Trans-1,2-Dichloroethene	ND	0.252	ND	1.00	12/13/18	KCA	1
trans-1,3-Dichloropropene	ND	0.221	ND	1.00	12/13/18	KCA	1
Trichloroethene	0.065	0.037	0.35	0.20	12/13/18	KCA	1
Trichlorofluoromethane	0.329	0.178	1.85	1.00	12/13/18	KCA	1
Trichlorotrifluoroethane	ND	0.131	ND	1.00	12/13/18	KCA	1
Vinyl Chloride	ND	0.078	ND	0.20	12/13/18	KCA	1
<u>QA/QC Surrogates</u>							
% Bromofluorobenzene	79	%	79	%	12/13/18	KCA	1

Client ID: SS 1

Parameter	ppbv Result	ppbv RL	ug/m3 Result	ug/m3 RL	Date/Time	By	Dilution
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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:

S - Laboratory solvent, contamination is possible.

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

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Phyllis Shiller, Laboratory Director

December 14, 2018

Reviewed and Released by: Rashmi Makol, Project Manager



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



QA/QC Report

December 14, 2018

QA/QC Data

SDG I.D.: GCC13973

Parameter	Blk ppbv	Blk RL ppbv	Blk ug/m3	Blk RL ug/m3	LCS %	Sample Result ug/m3	Sample Dup ug/m3	Sample Result ppbv	Sample Dup ppbv	DUP RPD	% Rec Limits	% RPD Limits
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QA/QC Batch 459950 (ppbv), QC Sample No: CC13975 (CC13973 (1X, 5X) , CC13974 (1X, 5X) , CC13975 (1X, 5X))

Volatiles

1,1,1,2-Tetrachloroethane	ND	0.150	ND	1.03	97	ND	ND	ND	ND	NC	70 - 130	25
1,1,1-Trichloroethane	ND	0.180	ND	0.98	85	ND	ND	ND	ND	NC	70 - 130	25
1,1,2,2-Tetrachloroethane	ND	0.150	ND	1.03	95	ND	ND	ND	ND	NC	70 - 130	25
1,1,2-Trichloroethane	ND	0.180	ND	0.98	97	ND	ND	ND	ND	NC	70 - 130	25
1,1-Dichloroethane	ND	0.250	ND	1.01	88	ND	ND	ND	ND	NC	70 - 130	25
1,1-Dichloroethene	ND	0.050	ND	0.20	83	ND	ND	ND	ND	NC	70 - 130	25
1,2,4-Trichlorobenzene	ND	0.130	ND	0.96	149	ND	ND	ND	ND	NC	70 - 130	25
1,2,4-Trimethylbenzene	ND	0.200	ND	0.98	98	4.81	4.62	0.979	0.941	NC	70 - 130	25
1,2-Dibromoethane(EDB)	ND	0.130	ND	1.00	98	ND	ND	ND	ND	NC	70 - 130	25
1,2-Dichlorobenzene	ND	0.170	ND	1.02	102	ND	ND	ND	ND	NC	70 - 130	25
1,2-Dichloroethane	ND	0.250	ND	1.01	85	ND	ND	ND	ND	NC	70 - 130	25
1,2-dichloropropane	ND	0.220	ND	1.02	95	ND	ND	ND	ND	NC	70 - 130	25
1,2-Dichlorotetrafluoroethane	ND	0.140	ND	0.98	92	ND	ND	ND	ND	NC	70 - 130	25
1,3,5-Trimethylbenzene	ND	0.200	ND	0.98	100	1.72	1.75	0.350	0.357	NC	70 - 130	25
1,3-Butadiene	ND	0.450	ND	0.99	90	ND	ND	ND	ND	NC	70 - 130	25
1,3-Dichlorobenzene	ND	0.170	ND	1.02	102	ND	ND	ND	ND	NC	70 - 130	25
1,4-Dichlorobenzene	ND	0.170	ND	1.02	100	ND	ND	ND	ND	NC	70 - 130	25
1,4-Dioxane	ND	0.280	ND	1.01	100	ND	ND	ND	ND	NC	70 - 130	25
2-Hexanone(MBK)	ND	0.240	ND	0.98	94	ND	ND	ND	ND	NC	70 - 130	25
4-Ethyltoluene	ND	0.200	ND	0.98	101	6.14	5.85	1.25	1.19	4.9	70 - 130	25
4-Isopropyltoluene	ND	0.180	ND	0.99	103	ND	ND	ND	ND	NC	70 - 130	25
4-Methyl-2-pentanone(MIBK)	ND	0.240	ND	0.98	91	1.44	1.54	0.352	0.377	NC	70 - 130	25
Acetone	ND	0.420	ND	1.00	82	131	129	55.1	54.5	1.1	70 - 130	25
Acrylonitrile	ND	0.460	ND	1.00	89	ND	ND	ND	ND	NC	70 - 130	25
Benzene	ND	0.310	ND	0.99	90	5.87	5.59	1.84	1.75	5.0	70 - 130	25
Benzyl chloride	ND	0.190	ND	0.98	106	ND	ND	ND	ND	NC	70 - 130	25
Bromodichloromethane	ND	0.150	ND	1.00	102	ND	ND	ND	ND	NC	70 - 130	25
Bromoform	ND	0.097	ND	1.00	103	ND	ND	ND	ND	NC	70 - 130	25
Bromomethane	ND	0.260	ND	1.01	86	ND	ND	ND	ND	NC	70 - 130	25
Carbon Disulfide	ND	0.320	ND	1.00	83	3.64	3.36	1.17	1.08	NC	70 - 130	25
Carbon Tetrachloride	ND	0.032	ND	0.20	88	0.33	0.30	0.053	0.047	NC	70 - 130	25
Chlorobenzene	ND	0.220	ND	1.01	103	ND	ND	ND	ND	NC	70 - 130	25
Chloroethane	ND	0.380	ND	1.00	85	ND	ND	ND	ND	NC	70 - 130	25
Chloroform	ND	0.200	ND	0.98	89	ND	ND	ND	ND	NC	70 - 130	25
Chloromethane	ND	0.480	ND	0.99	82	ND	ND	ND	ND	NC	70 - 130	25
Cis-1,2-Dichloroethene	ND	0.050	ND	0.20	89	ND	ND	ND	ND	NC	70 - 130	25
cis-1,3-Dichloropropene	ND	0.220	ND	1.00	98	ND	ND	ND	ND	NC	70 - 130	25
Cyclohexane	ND	0.290	ND	1.00	92	1.07	1.06	0.310	0.309	NC	70 - 130	25
Dibromochloromethane	ND	0.120	ND	1.02	98	ND	ND	ND	ND	NC	70 - 130	25
Dichlorodifluoromethane	ND	0.200	ND	0.99	93	1.50	1.45	0.304	0.293	NC	70 - 130	25
Ethanol	ND	0.530	ND	1.00	99	19.8	17.5	10.5	9.27	12.4	70 - 130	25

QA/QC Data

SDG I.D.: GCC13973

Parameter	Bik ppbv	Bik RL ppbv	Bik ug/m3	Bik RL ug/m3	LCS %	Sample Result ug/m3	Sample Dup ug/m3	Sample Result ppbv	Sample Dup ppbv	DUP RPD	% Rec Limits	% RPD Limits
Ethyl acetate	ND	0.280	ND	1.01	80	ND	ND	ND	ND	NC	70 - 130	25
Ethylbenzene	ND	0.230	ND	1.00	101	5.47	5.64	1.26	1.30	3.1	70 - 130	25
Heptane	ND	0.240	ND	0.98	89	3.26	2.93	0.797	0.715	NC	70 - 130	25
Hexachlorobutadiene	ND	0.094	ND	1.00	123	ND	ND	ND	ND	NC	70 - 130	25
Hexane	ND	0.280	ND	0.99	89	1.32 S	ND	0.375 S	ND	NC	70 - 130	25
Isopropylalcohol	ND	0.410	ND	1.01	82	2.87	2.92	1.17	1.19	NC	70 - 130	25
Isopropylbenzene	ND	0.200	ND	0.98	100	ND	ND	ND	ND	NC	70 - 130	25
m,p-Xylene	ND	0.230	ND	1.00	99	18.3	18.4	4.22	4.25	0.7	70 - 130	25
Methyl Ethyl Ketone	ND	0.340	ND	1.00	78	12.5	12.9	4.24	4.38	3.2	70 - 130	25
Methyl tert-butyl ether(MTBE)	ND	0.280	ND	1.01	87	ND	ND	ND	ND	NC	70 - 130	25
Methylene Chloride	ND	0.860	ND	2.99	81	ND	ND	ND	ND	NC	70 - 130	25
n-Butylbenzene	ND	0.180	ND	0.99	99	1.22	1.26	0.223	0.229	NC	70 - 130	25
o-Xylene	ND	0.230	ND	1.00	104	6.47	6.29	1.49	1.45	2.7	70 - 130	25
Propylene	ND	0.580	ND	1.00	95	2.79	2.89	1.62	1.68	NC	70 - 130	25
sec-Butylbenzene	ND	0.180	ND	0.99	96	ND	ND	ND	ND	NC	70 - 130	25
Styrene	ND	0.230	ND	0.98	104	2.88	3.39	0.676	0.797	NC	70 - 130	25
Tetrachloroethene	ND	0.037	ND	0.25	104	5.73	6.14	0.846	0.906	6.8	70 - 130	25
Tetrahydrofuran	ND	0.340	ND	1.00	88	ND	ND	ND	ND	NC	70 - 130	25
Toluene	ND	0.270	ND	1.02	97	21.6	21.4	5.74	5.69	0.9	70 - 130	25
Trans-1,2-Dichloroethene	ND	0.250	ND	0.99	85	ND	ND	ND	ND	NC	70 - 130	25
trans-1,3-Dichloropropene	ND	0.220	ND	1.00	95	ND	ND	ND	ND	NC	70 - 130	25
Trichloroethene	ND	0.037	ND	0.20	100	0.35	0.30	0.065	0.056	NC	70 - 130	25
Trichlorofluoromethane	ND	0.180	ND	1.01	85	1.85	1.56	0.329	0.278	NC	70 - 130	25
Trichlorotrifluoroethane	ND	0.130	ND	1.00	86	ND	ND	ND	ND	NC	70 - 130	25
Vinyl Chloride	ND	0.078	ND	0.20	91	ND	ND	ND	ND	NC	70 - 130	25
% Bromofluorobenzene	98		98		105	79	83	79	83	NC	70 - 130	25

I = This parameter is outside laboratory LCS/LCSD 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
December 14, 2018

Friday, December 14, 2018

Criteria: None

State: NY

Sample Criteria Exceedances Report

GCC13973 - EBC

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	RL Criteria	Analysis Units
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*** No Data to Display ***

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
Tel. (860) 645-1102 Fax (860) 645-0823



Analysis Comments

December 14, 2018

SDG I.D.: GCC13973

The following analysis comments are made regarding exceptions to criteria not already noted in the Analysis Report or QA/QC Report: None.



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**CHAIN OF CUSTODY RECORD
 AIR ANALYSES**

800-827-5426
 email: greg@phoenixlabs.com

P.O. # _____ Page 1 of 1

Data Delivery: _____
 Fax #: _____
 Email: F.1.e
 Phone #: _____

Report to: Keith Butler
 Customer: EBC
 Address: _____

Invoice to: EBC
 Project Name: 188 East 135th Street, Bronx
 Requested Deliverable: RCP ASP CAT B
 MCP NJ Deliverables
 State where samples collected: NY

Phoenix ID #	Client Sample ID	THIS SECTION FOR LAB USE ONLY										Ambient/Indoor Air	Soil Gas	Grab (G) Composite (C)	TO-14	TO-15			
		Canister ID #	Canister Size (L)	Outgoing Canister Pressure ("Hg)	Incoming Canister Pressure ("Hg)	Flow Regulator ID #	Flow Controller Setting (ml/min)	Sampling Start Time	Sampling End Time	Sample Start Date	Canister Pressure at Start ("Hg)						Canister Pressure at End ("Hg)		
	Pid not use	23346	6.0	-30		4985	43												
139173	SV1	19589			-6	4959						10:08	11:56	12-12-18	-28	-5		X	
139174	SV2	11290			-3	5652						10:21	12:32	12-12-18	-30	-5		X	
139175	SS1	19425			-7	5693						10:15	12:35	12-12-18	-29	-5		X	

Relinquished by: David Rukk
 Accepted by: Greg Butler
 Date: 12-13-18
 Time: 11:30

Requested Criteria: _____
 Turnaround Time: _____
 24 Hour 48 F
 I attest that all media released received in good working condition of this document.

Quote Number: _____
 Signature: f