

July 13, 2020

US Environmental Protection Agency Office of Ecosystem Protection 5 Post Office Square – Suite 100 Boston, MA 02109-3912 Attention Ms. Shauna Little

Notice of Intent for Remediation General Permit

Re: Project: Wheeler Self Storage Facility

Site: 0 Wheeler Road

Burlington, Massachusetts

BL Project No.: 16D3253

Dear Ms. Little,

BL Companies is submitting this Notice of Intent (NOI) on behalf of Wheeler Self Storage LLC for a Remediation General Permit (RGP) under EPA's National Pollutant Discharge Elimination System (NPDES) program. The RGP is required to discharge groundwater containing regulated compounds encountered during construction dewatering activities for development of a new self-storage facility at the referenced property (hereafter the "Site").

This application is organized as follows:

- Cover Letter
- Figures
- Attachment A RGP NOI Application Form
- Attachment B Stream Stats Flow Statistics Report
- Attachment C MACRIS Report
- Attachment D Endangered and Critical Species Report
- Attachment E Effluent Data Estimations
- Attachment F Groundwater Data
- Attachment G Treatment System Specifications and Data Sheets

The Site is located in Burlington, Massachusetts, is currently undeveloped land, and is known as 0 Wheeler Road. The Site consists of two parcels of land that are separated by a private road known as Rounder Way (a.k.a. Commonwealth Executive Drive). The Site is bounded to the south by Wheeler Road, to the east by Blanchard Road (a paper road) and commercial/industrial property, to the west by undeveloped land and commercial property, and to the north by Interstate 95/Route 128 and commercial property.

A Site Location Map is included as Figure 1 and a property survey map is included as Figure 2.

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The Site totals approximately 11.03 acres of land and the two parcels are identified as Tax Parcel 52-4-0 and 52-6-0. No structures or impervious pavement are currently present on any portion of the Site, and there is currently no stormwater management or treatment being utilized in the existing condition. The pending construction activities will only take place on approximately two acres located at the eastern end of Tax Parcel 52-4-0. The rest of this parcel and all of Tax Parcel 52-6-0 will be preserved in its current condition via a Conservation Restriction granted to the Town of Burlington Conservation Commission. The limits of the Conservation Restriction are shown on Figure 2 and the Limit of Disturbance for construction activities is shown on Figure 3.

This Conservation Restriction is required by the following permits: Site Plan issued by the Planning Board and filed with the Town Clerk on May 25, 2018; Special Permit filed with the Town Clerk on May 25, 2018 and recorded at the Middlesex South Registry of Deeds as Book 71169 Page 489; and Superseding Order of Conditions issued by the Massachusetts Department of Environmental Protection (MADEP) and recorded at the Middlesex South Registry of Deeds as Book 73760 Page 445. The referenced decisions permit the development of an 119,651+/- square foot, five-story self-storage facility with associated surface parking and on-site infrastructure improvements.

The construction dewatering is required in order to install the foundation system for the proposed self-storage building at the east end of the Site. Groundwater in the vicinity of the proposed building is contaminated from an off-Site release(s), necessitating management and treatment of the dewatering discharge under the RGP.

Several phases of soil and groundwater investigation have been completed in anticipation of Site acquisition and development. The locations of soil borings, test pits, and monitoring wells completed at the Site are shown on the Site Investigation Plan for the development area provided as Figure 4. Regulated compounds have been detected in both soil and groundwater above MADEP reportable concentrations. The Site will be reported as a disposal Site under the Massachusetts Contingency Plan (MCP) by the applicant upon acquisition of the Site, which is scheduled to occur in the next week. Known contaminants in the groundwater at the Site include Chlorinated Volatile Organic Compounds (CVOCs), and several metals. The CVOCs are attributed to an upgradient release of CVOCs from the release(s) at the former High Voltage Energy Corporation/Bellofram facility (RTN 3-0000981), and possibly other off-site disposal sites. Wheeler Self Storage, LLC will file for Downgradient Property Status under the MCP due to the impact to the Site from the upgradient release(s) of CVOCs. A source of CVOCs has not been identified on the Site and there is no history of such usage. The metals concentrations in groundwater are attributed to naturally occurring conditions.

Summary of Soil and Groundwater Investigations

During previous investigations at the Site completed in both 2017 and 2020, BL Companies identified regulated compounds in both soil and groundwater. EPH compounds, aromatic VOCs, 2-Methylnapthalene, and metals have been detected in soils at the southeastern corner of the Site, at concentrations above and below reportable concentrations under the MCP. The presence of these

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compounds is believed to be related to the storage of construction equipment and/or placement of fill on the southeastern corner of the Site. These impacts have been noted to be above the water table, with minor impacts to groundwater in the form of C9-C12 Aliphatic hydrocarbons, and Naphthalene. During the construction of the proposed self-storage facility, the remediation of the soils containing regulated compounds above RCS-1 reportable concentrations will be completed via Release Abatement Measures under the MCP. Soil conditions are summarized as follows:

- Soil samples from SB-101 (0.5-1.6') and SB-101 (5-8.9') were collected and analyzed, as well as soil samples collected from the test pits TP-1, TP-2, TP-3, TP-4, TP-6, TP-7, TP-8 and TP-9. During the previous Phase II Investigation, SB-3 was identified within this area with low concentrations of MADEP VPH and elevated concentrations of MADEP EPH compounds. These previous VPH and EPH detections however were identified below the MADEP Reportable Concentrations in Soil (RCS-1) Criteria which applies to the Site.
- Low concentrations of EPH and VPH were identified in SB-101 (0.5-1.6'), above the observed groundwater, which were consistent with the findings of the previous Phase II Investigation sample from SB-3. These concentrations were identified as being well below the MADEP RCS-1 Criteria. The sample TP-1 (3-4') was identified with EPH compounds, including C19-C36 Aliphatic Hydrocarbons above the MassDEP RCS-1 Criteria. C9-C10-Aromatic Hydrocarbons, C9-C12, Aliphatic Hydrocarbons, Ethyl Benzene, Xylenes, Naphthalene, and C9-C18 Aliphatic Hydrocarbons were identified below the MassDEP RCS-1 Criteria. Additionally, in TP-7 and TP-8, C9-C10-Aromatic Hydrocarbons, C9-C12, Aliphatic Hydrocarbons, Ethyl Benzene, Xylenes, Naphthalene, and C9-C18 Aliphatic Hydrocarbons were identified below the MassDEP RCS-1 Criteria. Aromatic VOCs were identified in the sample SB-101 (0.5-1.6'), including 1,2,4-Trimethylbenzene, Xylenes, and Naphthalene, however the concentrations were below the RCS-1 Criteria, and total VOCs were less than 2 mg/kg. Only Naphthalene was detected for VOCs in the deeper SB-101 (5-8.9') sample, below the water, and below the MADEP RCS-1 Criteria.
- Additional samples from TP-2, TP-3, TP-4, TP-6, TP-9 and a shallow sample from TP-7 did not identify VOCs, VPH, EPH, or SVOCs above the laboratory detection limit.
- SVOCs were identified in the sample SB-101 (0.5-1.6'), including 2-Methylnapthalene, Bis(2-ethylexyl) phthalate, and Pyrene. 2-Methylnapthalene was detected at concentrations above the RCS-1 Criteria. In TP-1 (3-4') 2-Methylnapthalene was identified above the MassDEP RCS-1 Criteria. Acenaphthylene, Bis(2-ethylhexyl) phthalate, Fluoranthene, Fluorene, Naphthalene, Phenanthrene, and Pyrene were identified below the MADEP RCS-1 Criteria.
- The analytical results from these samples identified concentrations of Arsenic, Barium, Cadmium, Chromium, Nickel, Vanadium, and Zinc, however at concentrations below the RCS-1 Standard. These concentrations appear to be indicative of background, natural soil conditions, and do not represent a potential release condition at the Site. Lead was detected at a concentration above the MADEP RCS-1 Criteria.

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The source of the CVOC impacts identified in the groundwater is believed to be an upgradient release(s) at the former High Voltage Energy Corporation/Bellofram facility (RTN 3-0000981). A Downgradient Property Status report (DPS) is forthcoming. The VOCs trichloroethene (TCE) and tetrachloroethene (PCE) were detected above applicable MCP Reportable Concentrations for Groundwater (RCGW-1) in the groundwater sample collected from monitoring wells MW-2, MW-101, MW-102, and MW-103, with cis 1,2-dichloroethane detected at the same locations, below the RCGW-1 criteria. 1,3,5-Trimethylbenzene was detected below the MCP reportable concentrations in sample MW-102, and MW-103. 1,2,4-Trimethylbenzene, 1,3,5-Trimethylbenzene, Xylenes, and Naphthalene were identified in the sample collected from MW-101, below the applicable RCGW-1 Criteria. VOCs were not detected at concentrations above their respective laboratory reporting limits in MW-104.

These VOC detections were similar to those previously identified in 2017. The CVOC plume generally extends across the eastern portion of the Site. An on-Site source of CVOCs has not been identified and the highest concentrations of CVOCs in groundwater are present at the upgradient property boundary with respect to groundwater flow. The presence of the CVOC plume will be reported under the MCP and the applicant intends to file for Downgradient Property Status. The proposed building will be constructed with a chemical vapor barrier and a sub-slab depressurization system to mitigate any potential vapor intrusion concerns.

Proposed Dewatering Discharge Treatment and Discharge

The dewatering discharge will be treated and subsequently discharged directly to an unnamed tributary to Vine Brook. The proposed discharge location is shown on Figure 2. Vine Brook is a Class B surface water body that flows north from the Site under I-95/Route 128 and through a highly developed commercial area. There is no alternative other than direct discharge of the dewatering effluent to Vine Brook, as the Town of Burlington will not allow discharge to the storm sewer or municipal sewer systems. Additional information regarding the receiving waters is included in Attachment B.

The most recent groundwater quality data was collected in March and May 2020 for Metals, VOCs, Extractable Petroleum Hydrocarbons (EPH), Volatile Petroleum Hydrocarbons (VPH), and Semi-Volatile Organic Compounds (SVOCs). Copies of the most recent laboratory analytical reports for groundwater in the vicinity of the proposed construction dewatering are included in Attachment F.

The proposed groundwater treatment system for this project consists of influent waters pumped from the immediate vicinity of construction excavations that will extend below the water table using a well-point system. The well-points will pump through a central manifold into an 18,000-gallon weir tank. The weir tank will be equipped with a pH adjustment system and chemical aided settling system to enhance particulate removal from the discharge. The contents of the tank will then be pumped through bag filtration to remove the metals-containing sediment, through two Granulated Activated Carbon (GAC) vessels for VOC removal, and through a Resin media vessel

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for final polishing and treatment. Effluent flows will be metered for total flow documentation. The treatment system design components are illustrated in Figure 5.

The treated groundwater will then be discharged to the nearby surface water body on the adjacent western parcel, which flows to the nearby Vine Brook, ultimately discharging to the wetlands to the north of the Site. The discharge location is shown on Figure 2.

Dewatering is intended to be intermittent during the construction as groundwater is encountered during footing and foundation installation. Additional information regarding the treatment system is included in Attachments E and G.

Endangered and Critical Species Evaluation

BL Companies reviewed information through the Massachusetts Division of Fisheries and Wildlife (MADFW) Massachusetts Priority Resource Map (Figure 6) and through the US Fish and Wildlife Service for information on Endangered and Critical Species. Based on the findings, the Site and the location of proposed discharge do not appear to be locations within an Area of Critical Environmental Concern.

The review of information through the MADFW and the US Fish and Wildlife Service concluded that the Site falls under Criterion C of the RGP guidance. The guidance for Criterion C includes the following:

- Criterion C: Discharges not likely to adversely affect ESA-listed species and/or designated critical habitat. ESA-listed species and/or designated critical habitat(s) under the jurisdiction of the USFWS and/or NMFS are likely to occur in or near your site's "action area," and you certify to EPA that your site's discharges and discharge-related activities are not likely to adversely affect ESA-listed threatened or endangered species and/or designated critical habitat. This certification may include consideration of any stormwater controls and/or management practices you will adopt to ensure that your discharges and discharge-related activities are not likely to adversely affect ESA-listed species and/or designated critical habitat. To certify your eligibility under this criterion, indicate 1) the ESA-listed species and/or designated habitat located in your "action area" using the process outlined in Attachment D of this permit; 2) the distance between the site and the listed species and/or designated critical habitat in the action area (in miles); and 3) a rationale describing specifically how adverse effects to ESA-listed species will be avoided from the discharges and discharge-related activities. You must also include a copy of your site map from your SWPPP showing the upland and in-water extent of your "action area" with this NOI.
- Basis statement content/Supporting documentation: A basis statement supporting the selection of Criterion C should identify the information resources and expertise (e.g., state or federal biologists) used to arrive at this conclusion. Any supporting documentation

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should explicitly state that both ESA-listed species and designated critical habitat under the jurisdiction of the USFWS and/or NMFS were considered in the evaluation.

In accordance with the guidance for Criterion C, the findings of the endangered and critical species evaluation are as follows:

Resources Consulted:

United States Fish and Wildlife Service Information for Planning and Consultation (IPaC) tool, Massachusetts Natural Heritage and Endangered Species Program online maps (https://www.mass.gov/service-details/the-northern-long-eared-bat), National Marine Fisheries Service (NMFS) Essential Fish Habitat map, and NMFS CGP Waters of Concern Map.

ESA-listed Species/Critical Habitat in Action Area:

One species, the Northern Long-Eared Bat (*Myotis septentrionalis*), or NLEB, was identified as potentially present in the vicinity of the action area according to the US Fish and Wildlife Service (USFWS) Information for Planning and Consultation (IPaC) response. The NLEB is a forest interior species that is considered to be a "clutter specialist" as they are capable of foraging in dense, interior forests. According to NatureServe.org, this bat generally is associated with old-growth forests composed of trees 100 years old or older. It relies on intact interior forest habitat, with low edge-to-interior ratios. Relevant late-successional forest features include a high percentage of old trees, uneven forest structure (resulting in multilayered vertical structure), single and multiple tree-fall gaps, standing snags, and woody debris. During the summer, the species is widely distributed in a variety of wooded habitats and its habitat needs for roosting and foraging appear to be flexible and not specific. Thus, the USFWS does not have any critical habitat mapped for NLEB. Therefore, there is no designated critical habitat for the NLEB present within the action area.

Furthermore, according to the USFWS critical habitat mapper, there is no other critical habitats mapped within the action area. The action area is located in a relatively developed commercial area, adjacent to a major highway, making it unlikely that this species is present within the vicinity of the action area. While there are mature trees present on-site, they have been identified primarily within the portions of the Site slated for Conservation Restriction. In addition, this forested habitat is not considered "intact interior forest" and does not consist of uneven forest structures, trees with single and/or multiple tree-fall gaps, or standing snags. The aforementioned habitat is located very close to roads, a highway, and commercial office buildings and parking lots which are likely not preferred by the NLEB due to the fragmented, forest-edge condition. The presence of non-native invasive plant species exacerbates these limitations, as they are not ideal habitat or food sources for native wildlife. Larger blocks of mature forest are present in an approximate 9-acre portion of the Site to the west of the action area that has been protected via a conservation easement with the Town of Burlington, in a conservation restriction along the southern side of Wheeler Road and in an even larger forested and park area that is present approximately 0.25 mile to the east and southeast, which would be more appropriate habitat for the NLEB.

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Distance Between Site and ESA-listed Species/Critical Habitat:

According to mapping from Massachusetts Natural Heritage and Endangered Species Program (NHESP), the nearest known NLEB hibernacula is are located approximately 7.3 miles northwest of the Site, and the nearest known maternity roost tree is located approximately 63 miles southeast of the Site.

How Adverse Effects will be Avoided:

Given the lack of appropriate interior forest habitat, it is highly unlikely that the NLEB is present on-site in its existing conditions. However, natural conditions on-site will be improved by the addition and enhancement of landscaped areas, and addition of stormwater treatments including a rain garden that will naturally filter stormwater before it is returned to the groundwater table or to the adjacent tributary to Vine Brook. The eastern portion of the Site where the proposed development will be located is highly disturbed and degraded due to historic land use for gravel mining and crane and construction equipment storage. Various refuse and bulky debris are noted throughout this area and consist of smaller construction equipment, concrete, typical domestic refuse, and mechanical refuse such as car batteries, and will be removed from the Site and revegetated. Non-native invasive plant species do limit the available habitat on-site as they do not provide food sources or shelter for native wildlife. The non-native plants in the eastern portion of the Site are overtaking native plant species, namely autumn olive (E. umbrellata), glossy buckhorn (F. alnus), and oriental bittersweet (C. orbiculatus). Additionally, the natural buffer along the tributary to Vine Brook will be enhanced with additional native plantings providing future habitat for roosting and foraging opportunities for this species. The addition of plantings and stormwater treatment will offer an improvement to the water quality being discharged from the Site into the Vine Brook tributary and may provide an improvement to water quality for NLEB, and other terrestrial and aquatic species that may utilize Vine Brook as a water source.

Cultural and Historic Resources Evaluation

BL Companies reviewed information through the Massachusetts Cultural Resource Information Viewer (MACRIS) regarding the potential presence of significant cultural resources in the vicinity of the Site. The MACRIS report is included in Attachment C. Based on the information reviewed, the Site and the location of proposed discharge do not appear to be within areas of historical significance through the MACRIS system.

Closing

The NOI Application is included as Attachment A. The information provided on the application form and in the attached information demonstrates that the proposed dewatering discharge is eligible for authorization under the RGP.

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Should you have any questions, do not hesitate to contact the undersigned.

Sincerely, BL Companies

Samuel R. Haydock, MS, LEP

Principal

CC: Cathy Vakalopoulos – MADEP

Samuel Weissman - Wheeler Self Storage, LLC

Attachments:

Figure 1 – Site Location Map

Figure 2 – Conservation Restriction and Easement Map

Figure 3 – Overall Limit of Disturbance Area

Figure 4 – Site Investigation Plan

 $Figure \ 5-Treatment \ System \ Design$

Figure 6 – MADFW Priority Resource Map

Attachment A – RGP NOI Application Form

Attachment B – Stream Stats Flow Statistics Report

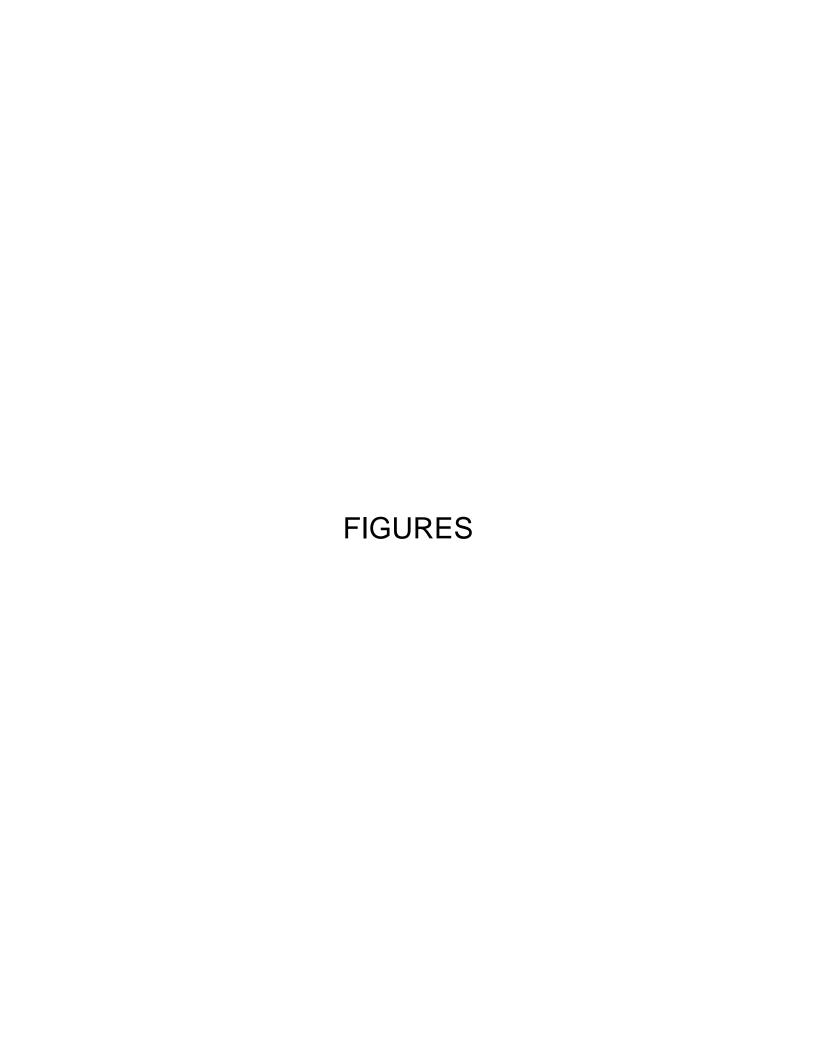
Attachment C – MACRIS Report

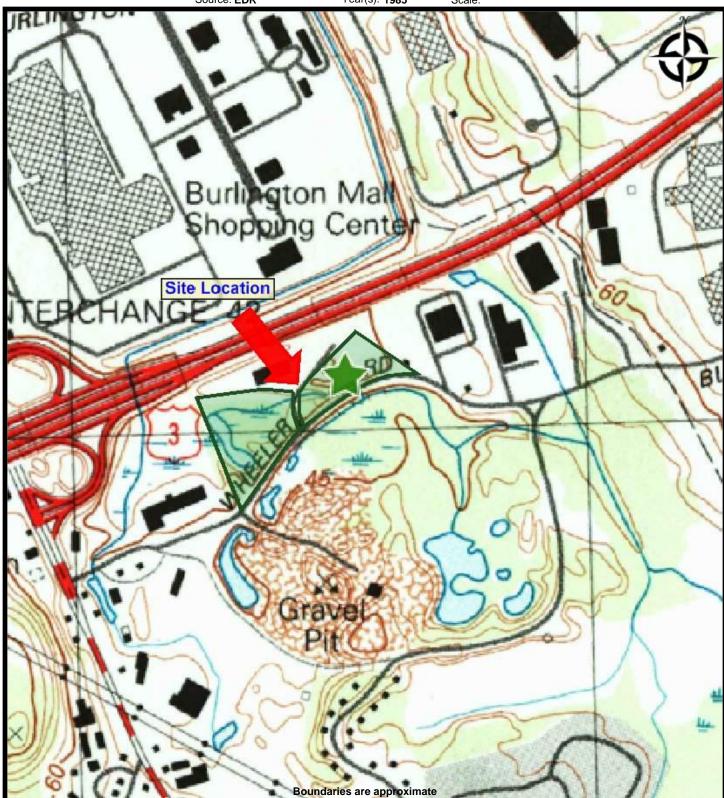
Attachment D – Endangered and Critical Species Report

 $Attachment \ E-Effluent \ Data \ Estimations$

Attachment F - Groundwater Data

 $\label{lem:system} \textbf{Attachment } \textbf{G}-\textbf{Treatment System Specifications and Data Sheets}$



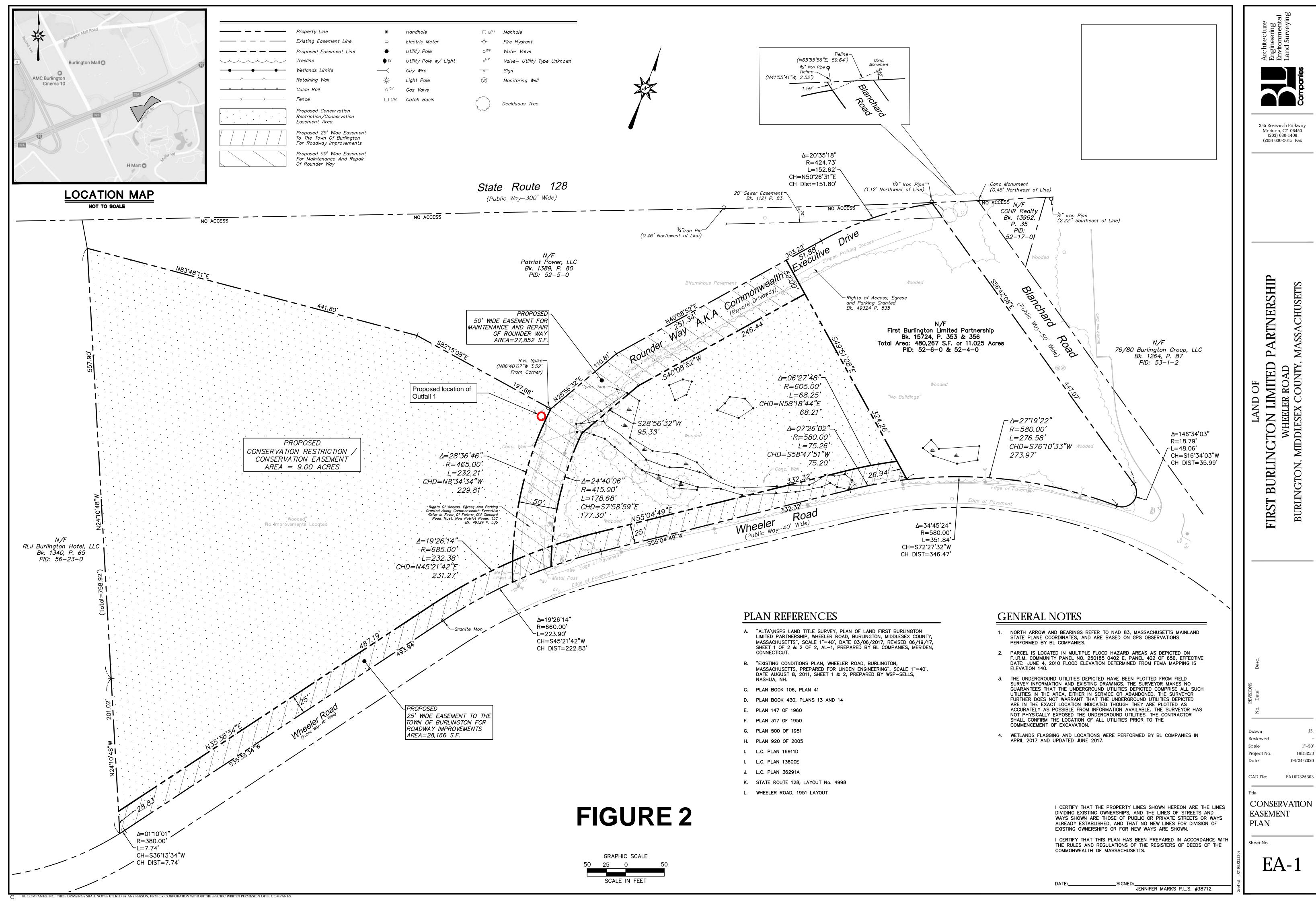




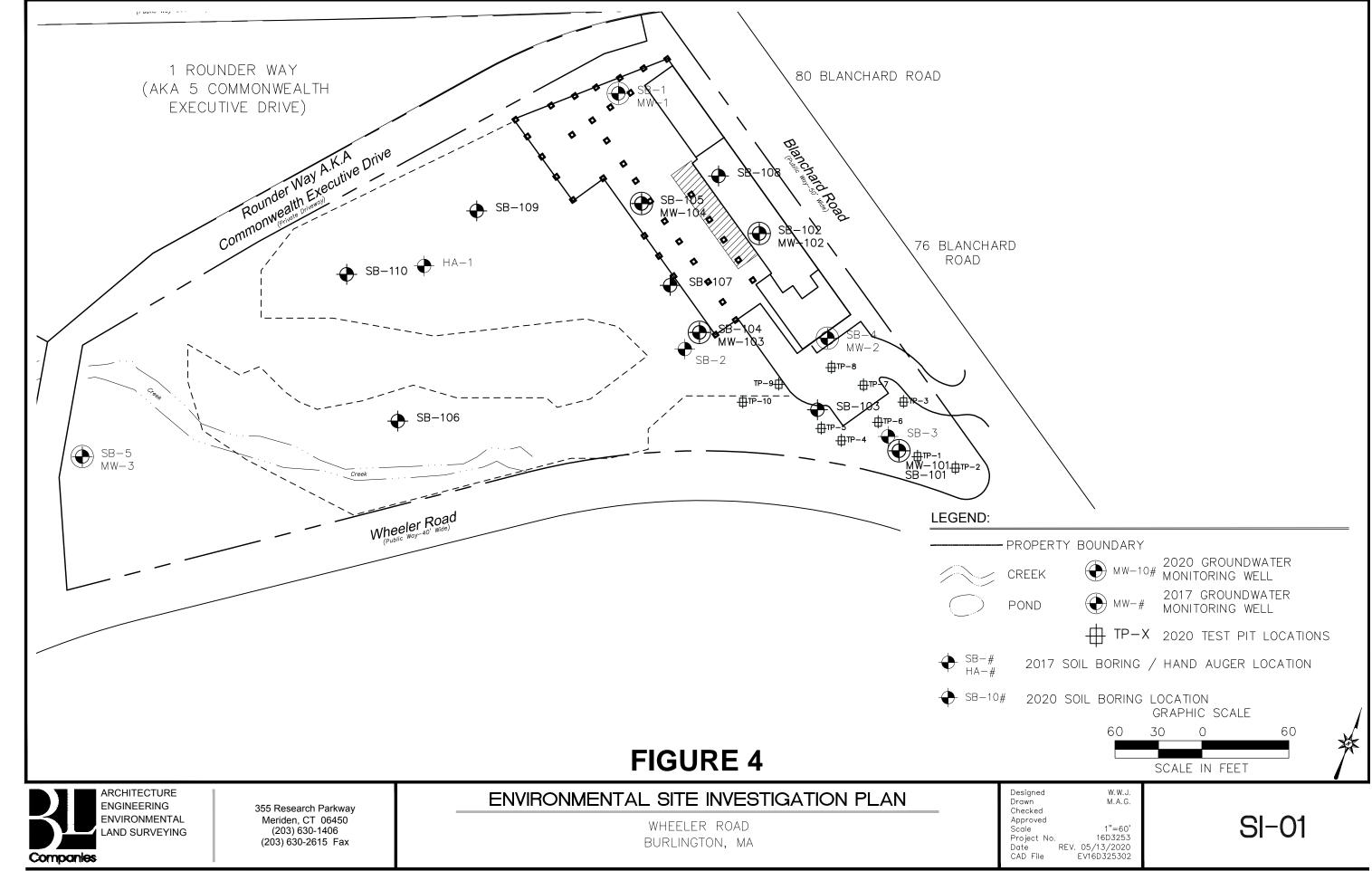
Architecture Engineering Environmental Land Surveying FIGURE 1 - SITE LOCATION MAP
SELF-STORAGE FACILTY SITE
Wheeler Road
Burlington, Massachusetts 01803

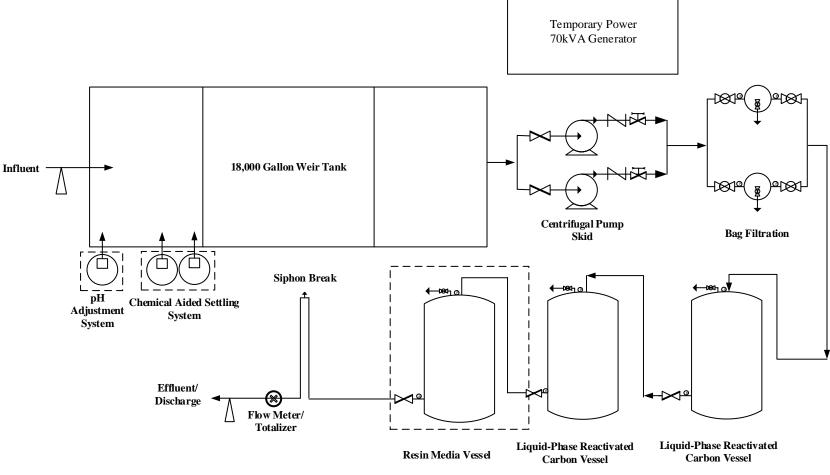
IMAGE YEAR: 1985

PROJ. #: 16D3253





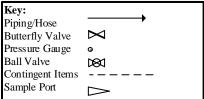




Notes:

- 1.) Figure is not to scale
- 2.) System rated for 250-300 GPM

FIGURE 5





Lockwood Remediation Technologies, LLC 89 Crawford Street Leominster, MA 01453

Leominster, MA 01453 Office: 774-450-7177

DESIGNED BY: LRT DRAWN BY: JHJ

CHECKED BY: DATE:

Water Treatment System Schematic

Public Storage Building Wheeler Road Burlington, MA

3224

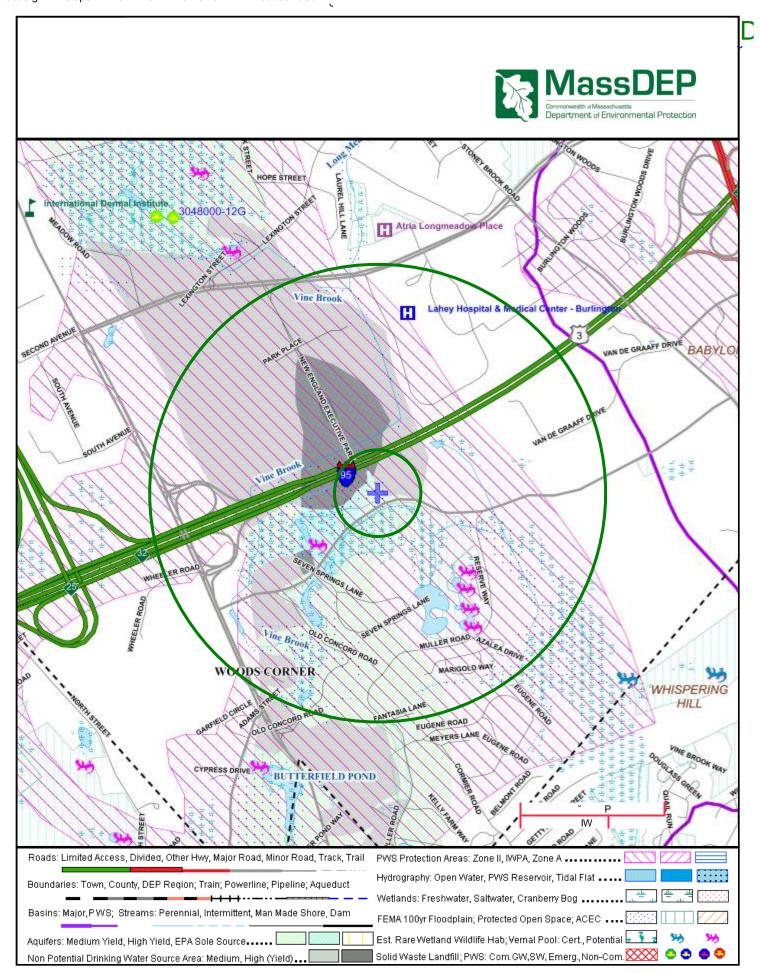


FIGURE 6

ATTACHMENT A NOI FORM

II. Suggested Format for the Remediation General Permit Notice of Intent (NOI)

A. General site information:

1. Name of site:	Site address:			
	Street:			
	City:		State:	Zip:
2. Site owner	Contact Person:			
	Telephone:	Email:		
	Mailing address:			
	Street:			
Owner is (check one): ☐ Federal ☐ State/Tribal ☐ Private ☐ Other; if so, specify:	City:		State:	Zip:
3. Site operator, if different than owner	Contact Person:			
	Telephone:	Email:		
	Mailing address:			
	Street:			
	City:		State:	Zip:
4. NPDES permit number assigned by EPA:	5. Other regulatory program(s) that apply to the site	(check all th	at apply):	
	☐ MA Chapter 21e; list RTN(s):	□ CERCL	.A	
NPDES permit is (check all that apply: \square RGP \square DGP \square CGP	☐ NH Groundwater Management Permit or	□ UIC Pro	•	
☐ MSGP ☐ Individual NPDES permit ☐ Other; if so, specify:	Groundwater Release Detection Permit:		Pretreatment	t
· · · · · · · · · · · · · · · · · · ·		□ CWA S	ection 404	

B	Receiving water information:
1	Name of receiving water(s).

1. Name of receiving water(s):	Waterbody identification of receiving water	(s): Classific	ation of receiving water(s):
Receiving water is (check any that apply): □ Outstar	nding Resource Water □ Ocean Sanctuary □ territo	rial sea □ Wild and Scenic Ri	ver
2. Has the operator attached a location map in accord	lance with the instructions in B, above? (check one)	: □ Yes □ No	
Are sensitive receptors present near the site? (check of the sensitive receptors) that is the sensitive receptors present near the site? (check of the sensitive receptors) are sensitive receptors present near the site?	one): □ Yes □ No		
3. Indicate if the receiving water(s) is listed in the Stapollutants indicated. Also, indicate if a final TMDL in 4.6 of the RGP.			
4. Indicate the seven day-ten-year low flow (7Q10) of Appendix V for sites located in Massachusetts and A		n the instructions in	
5. Indicate the requested dilution factor for the calculaccordance with the instructions in Appendix V for s			
6. Has the operator received confirmation from the a If yes, indicate date confirmation received:7. Has the operator attached a summary of receiving	-		
(check one): ☐ Yes ☐ No			
C. Source water information:			
1. Source water(s) is (check any that apply):			
☐ Contaminated groundwater	☐ Contaminated surface water	☐ The receiving water	☐ Potable water; if so, indicate municipality or origin:
Has the operator attached a summary of influent sampling results as required in Part 4.2 of the RGP	Has the operator attached a summary of influent sampling results as required in Part 4.2 of the	☐ A surface water other	
in accordance with the instruction in Appendix VIII? (check one):	RGP in accordance with the instruction in Appendix VIII? (check one):	than the receiving water; if so, indicate waterbody:	☐ Other; if so, specify:
□ Yes □ No	□ Yes □ No		

2. Source water contaminants:	
a. For source waters that are contaminated groundwater or contaminated surface water, indicate are any contaminants present that are not included in	b. For a source water that is a surface water other than the receiving water, potable water or other, indicate any contaminants present at the maximum concentration in accordance
the RGP? (check one): ☐ Yes ☐ No If yes, indicate the contaminant(s) and the maximum concentration present in accordance with the instructions in Appendix VIII.	with the instructions in Appendix VIII? (check one): □ Yes □ No
3. Has the source water been previously chlorinated or otherwise contains resid	dual chlorine? (check one): ☐ Yes ☐ No
D. Discharge information	
1.The discharge(s) is a(n) (check any that apply): \Box Existing discharge \Box New	w discharge □ New source
Outfall(s):	Outfall location(s): (Latitude, Longitude)
Discharges enter the receiving water(s) via (check any that apply): □ Direct di	scharge to the receiving water \Box Indirect discharge, if so, specify:
☐ A private storm sewer system ☐ A municipal storm sewer system If the discharge enters the receiving water via a private or municipal storm sew	ver system:
Has notification been provided to the owner of this system? (check one): □ Ye	•
Has the operator has received permission from the owner to use such system for obtaining permission:	or discharges? (check one): \square Yes \square No, if so, explain, with an estimated timeframe for
Has the operator attached a summary of any additional requirements the owner	of this system has specified? (check one): \square Yes \square No
Provide the expected start and end dates of discharge(s) (month/year):	
Indicate if the discharge is expected to occur over a duration of: \Box less than 1	2 months □ 12 months or more □ is an emergency discharge
Has the operator attached a site plan in accordance with the instructions in D, a	above? (check one): Yes No

2. Activity Category: (check all that apply)	3. Contamination Type Category: (check	all that apply)
	a. If Activity Categ	ory I or II: (check all that apply)
	 □ A. Inorganics □ B. Non-Halogenated Volatile Organic □ C. Halogenated Volatile Organic Cor □ D. Non-Halogenated Semi-Volatile Organic □ E. Halogenated Semi-Volatile Organi □ F. Fuels Parameters 	mpounds Organic Compounds
 □ I – Petroleum-Related Site Remediation □ II – Non-Petroleum-Related Site Remediation 	b. If Activity Category III, IV	V, V, VI, VII or VIII: (check either G or H)
 □ III – Non-Petroleum-Related Site Remediation □ III – Contaminated Site Dewatering □ IV – Dewatering of Pipelines and Tanks □ V – Aquifer Pump Testing □ VI – Well Development/Rehabilitation □ VII – Collection Structure Dewatering/Remediation □ VIII – Dredge-Related Dewatering 	□ G. Sites with Known Contamination c. If Category III-G, IV-G, V-G, VI-G, VII-G or VIII-G: (check all that apply) □ A. Inorganics □ B. Non-Halogenated Volatile Organic Compounds □ C. Halogenated Volatile Organic Compounds □ D. Non-Halogenated Semi-Volatile Organic Compounds □ E. Halogenated Semi-Volatile Organic Compounds □ F. Fuels Parameters	□ H. Sites with Unknown Contamination d. If Category III-H, IV-H, V-H, VI-H, VII-H or VIII-H Contamination Type Categories A through F apply

4. Influent and Effluent Characteristics

	Known	Known		75 5 4	5	Infl	uent	Effluent Lir	nitations
Parameter	or believed absent	or believed present	# of samples	Test method (#)	Detection limit (µg/l)	Daily maximum (µg/l)	Daily average (µg/l)	TBEL	WQBEL
A. Inorganics									
Ammonia								Report mg/L	
Chloride								Report µg/l	
Total Residual Chlorine								0.2 mg/L	
Total Suspended Solids								30 mg/L	
Antimony								206 μg/L	
Arsenic								104 μg/L	
Cadmium								10.2 μg/L	
Chromium III								323 μg/L	
Chromium VI								323 μg/L	
Copper								242 μg/L	
Iron								5,000 μg/L	
Lead								160 μg/L	
Mercury								0.739 μg/L	
Nickel								1,450 μg/L	
Selenium								235.8 μg/L	
Silver								35.1 μg/L	
Zinc								420 μg/L	
Cyanide								178 mg/L	
B. Non-Halogenated VOCs	3								
Total BTEX								100 μg/L	
Benzene								5.0 μg/L	
1,4 Dioxane								200 μg/L	
Acetone								7.97 mg/L	
Phenol								1,080 µg/L	

	Known	Known		_	_	Inf	luent	Effluent Lin	nitations
Parameter	or believed absent	or believed present	# of samples	Test method (#)	Detection limit (µg/l)	Daily maximum (µg/l)	Daily average (µg/l)	TBEL	WQBEL
C. Halogenated VOCs									
Carbon Tetrachloride								4.4 μg/L	
1,2 Dichlorobenzene								600 μg/L	
1,3 Dichlorobenzene								320 μg/L	
1,4 Dichlorobenzene								5.0 μg/L	
Total dichlorobenzene								763 µg/L in NH	
1,1 Dichloroethane								70 μg/L	
1,2 Dichloroethane								5.0 μg/L	
1,1 Dichloroethylene								3.2 µg/L	
Ethylene Dibromide								0.05 μg/L	
Methylene Chloride								4.6 μg/L	
1,1,1 Trichloroethane								200 μg/L	
1,1,2 Trichloroethane								5.0 μg/L	
Trichloroethylene								5.0 μg/L	
Tetrachloroethylene								5.0 μg/L	
cis-1,2 Dichloroethylene								70 μg/L	
Vinyl Chloride								2.0 μg/L	
D. Non-Halogenated SVO	Cs								
Total Phthalates								190 μg/L	
Diethylhexyl phthalate								101 μg/L	
Total Group I PAHs								1.0 μg/L	
Benzo(a)anthracene								_	
Benzo(a)pyrene								_	
Benzo(b)fluoranthene								_	
Benzo(k)fluoranthene								As Total PAHs	
Chrysene								_	
Dibenzo(a,h)anthracene								_	
Indeno(1,2,3-cd)pyrene									

	Known	Known				Inf	luent	Effluent Lin	nitations
Parameter	or believed absent	or believed present	# of samples	Test method (#)	Detection limit (µg/l)	Daily maximum (µg/l)	Daily average (µg/l)	TBEL	WQBEL
Total Group II PAHs								100 μg/L	
Naphthalene								20 μg/L	
E. Halogenated SVOCs									
Total PCBs								0.000064 µg/L	
Pentachlorophenol								1.0 μg/L	
	1			•					
F. Fuels Parameters Total Petroleum		1	1	1		1 1		<u> </u>	
Hydrocarbons								5.0 mg/L	
Ethanol								Report mg/L	
Methyl-tert-Butyl Ether								70 μg/L	
tert-Butyl Alcohol								120 μg/L in MA 40 μg/L in NH	
tert-Amyl Methyl Ether								90 μg/L in MA 140 μg/L in NH	
Other (i.e., pH, temperatur	re, hardness,	salinity, LC	50, addition	al pollutar	ats present);	if so, specify:			

E. Treatment system information

1. Indicate the type(s) of treatment that will be applied to effluent prior to discharge: (check all that apply)	
☐ Adsorption/Absorption ☐ Advanced Oxidation Processes ☐ Air Stripping ☐ Granulated Activated Carbon ("GAC")/Liquid Phase Carbon Adsorption	
☐ Ion Exchange ☐ Precipitation/Coagulation/Flocculation ☐ Separation/Filtration ☐ Other; if so, specify:	
2. Provide a written description of all treatment system(s) or processes that will be applied to the effluent prior to discharge.	
Identify each major treatment component (check any that apply):	
☐ Fractionation tanks☐ Equalization tank ☐ Oil/water separator ☐ Mechanical filter ☐ Media filter	
☐ Chemical feed tank ☐ Air stripping unit ☐ Bag filter ☐ Other; if so, specify:	
Indicate if either of the following will occur (check any that apply):	
□ Chlorination □ De-chlorination	
3. Provide the design flow capacity in gallons per minute (gpm) of the most limiting component.	
Indicate the most limiting component:	
Is use of a flow meter feasible? (check one): \square Yes \square No, if so, provide justification:	
Provide the proposed maximum effluent flow in gpm.	
Trovide the proposed maximum errident now in gpin.	
Provide the average effluent flow in gpm.	
If Activity Category IV applies, indicate the estimated total volume of water that will be discharged:	
4. Has the operator attached a schematic of flow in accordance with the instructions in E, above? (check one): ☐ Yes ☐ No	

F. Chemical and additive information

1. Indicate the type(s) of chemical or additive that will be applied to effluent prior to discharge or that may otherwise be present in the discharge(s): (check all that apply)
□ Algaecides/biocides □ Antifoams □ Coagulants □ Corrosion/scale inhibitors □ Disinfectants □ Flocculants □ Neutralizing agents □ Oxidants □ Oxygen □
scavengers □ pH conditioners □ Bioremedial agents, including microbes □ Chlorine or chemicals containing chlorine □ Other; if so, specify:
2. Provide the following information for each chemical/additive, using attachments, if necessary:
a. Product name, chemical formula, and manufacturer of the chemical/additive; b. Purpose or use of the chemical/additive or remedial agent; c. Material Safety Data Sheet (MSDS) and Chemical Abstracts Service (CAS) Registry number for each chemical/additive; d. The frequency (hourly, daily, etc.), duration (hours, days), quantity (maximum and average), and method of application for the chemical/additive; e. Any material compatibility risks for storage and/or use including the control measures used to minimize such risks; and f. If available, the vendor's reported aquatic toxicity (NOAEL and/or LC50 in percent for aquatic organism(s)).
3. Has the operator attached an explanation which demonstrates that the addition of such chemicals/additives may be authorized under this general permit in accordance
with the instructions in F, above? (check one): \square Yes \square No; if no, has the operator attached data that demonstrates each of the 126 priority pollutants in CWA Section 307(a) and 40 CFR Part 423.15(j)(1) are non-detect in discharges with the addition of the proposed chemical/additive?
(check one): ☐ Yes ☐ No
G. Endangered Species Act eligibility determination
1. Indicate under which criterion the discharge(s) is eligible for coverage under this general permit:
□ FWS Criterion A : No endangered or threatened species or critical habitat are in proximity to the discharges or related activities or come in contact with the "action area".
□ FWS Criterion B : Formal or informal consultation with the FWS under section 7 of the ESA resulted in either a no jeopardy opinion (formal consultation) or a written concurrence by FWS on a finding that the discharges and related activities are "not likely to adversely affect" listed species or critical habitat
(informal consultation). Has the operator completed consultation with FWS? (check one): ☐ Yes ☐ No; if no, is consultation underway? (check one): ☐
Yes □ No
□ FWS Criterion C : Using the best scientific and commercial data available, the effect of the discharges and related activities on listed species and critical habitat have been evaluated. Based on those evaluations, a determination is made by EPA, or by the operator and affirmed by EPA, that the discharges and related activities will have "no effect" on any federally threatened or endangered listed species or designated critical habitat under the jurisdiction of the FWS. This determination was made by: (check one) □ the operator □ EPA □ Other; if so, specify:
1 11.5. This determination was made by, (effect one) in the operator in the A in Other, it so, specify.

□ NMFS Criterion : A determination made by EPA is affirmed by the operator that the discharges and related activities will have "no effect" or are "not likely to adversely affect" any federally threatened or endangered listed species or critical habitat under the jurisdiction of NMFS and will not result in any take of
listed species. Has the operator previously completed consultation with NMFS? (check one): ☐ Yes ☐ No
2. Has the operator attached supporting documentation of ESA eligibility in accordance with the instructions in Appendix I, and G, above? (check one): \square Yes \square No
Does the supporting documentation include any written concurrence or finding provided by the Services? (check one): ☐ Yes ☐ No; if yes, attach.
H. National Historic Preservation Act eligibility determination
1. Indicate under which criterion the discharge(s) is eligible for coverage under this general permit:
□ Criterion A : No historic properties are present. The discharges and discharge-related activities (e.g., BMPs) do not have the potential to cause effects on historic properties.
☐ Criterion B: Historic properties are present. Discharges and discharge related activities do not have the potential to cause effects on historic properties.
□ Criterion C : Historic properties are present. The discharges and discharge-related activities have the potential to have an effect or will have an adverse effect on historic properties.
2. Has the operator attached supporting documentation of NHPA eligibility in accordance with the instructions in H, above? (check one): ☐ Yes ☐ No
Does the supporting documentation include any written agreement with the State Historic Preservation Officer (SHPO), Tribal Historic Preservation Officer (TPHO), or
other tribal representative that outlines measures the operator will carry out to mitigate or prevent any adverse effects on historic properties? (check one): \square Yes \square No
I. Supplemental information
Describe any supplemental information being provided with the NOI. Include attachments if required or otherwise necessary.
Has the operator attached data, including any laboratory case narrative and chain of custody used to support the application? (check one): ☐ Yes ☐ No
Has the operator attached the certification requirement for the Best Management Practices Plan (BMPP)? (check one): ☐ Yes ☐ No

J. Certification requirement

I certify under penalty of law that this document and all attachments were prepared under my direction or supervision in accordance with a system designed to assur that qualified personnel properly gathered and evaluated the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I have no personal knowledge that the information submitting than true, accurate, and complete. I am aware that there are significant penalties for submitting false information, including the possibility of fine and imprisonment for knowing violations.	ccordance with a syster versons who manage th lief, true, accurate, an significant penalties f	n designed to assur te system, or those d complete. I have or submitting false
A BMPP meeting the requirements of this general permit will be implemented prior to initiating BMPP certification statement: discharge.	mented prior to ir	nitiating
Notification provided to the appropriate State, including a copy of this NOI, if required.	Check one: Yes	No 🗆
Notification provided to the municipality in which the discharge is located, including a copy of this NOI, if requested.	Check one: Yes	No 🔳
Notification provided to the owner of a private or municipal storm sewer system, if such system is used for site discharges, including a copy of this NOI. if requested.	Check one: Yes □	No□ NA ■
Permission obtained from the owner of a private or municipal storm sewer system, if such system is used for site discharges. If yes, attach additional conditions. If no, attach explanation and timeframe for obtaining permission.	Check one: Yes □	No□ NA 🔳
Notification provided to the owner/operator of the area associated with activities covered by an additional discharge permit is (check one); □ RGP □ DGP □ CGP □ MSGP □ Individual NPDES nermit	Check one Ved	₩ VIV.
□ Other; if so, specify:		
Signature: My Date:	8/13/2020	
Print Name and Title: Samuel Weissman - Agent for Owner		

ATTACHMENT B STREAMSTATS REPORT

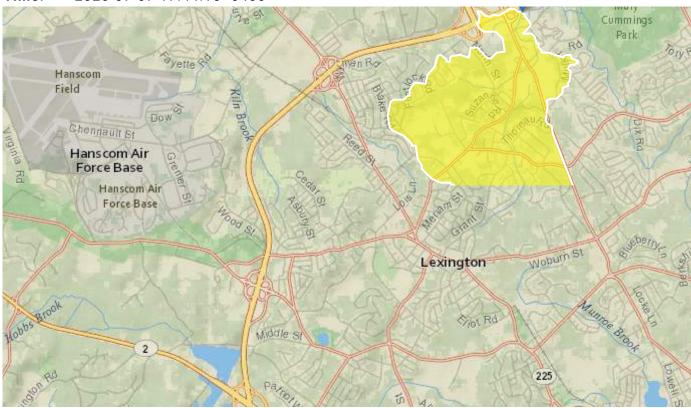
StreamStats Report

Region ID: MA

Workspace ID: MA20200707214402183000

Clicked Point (Latitude, Longitude): 42.47786, -71.21190

Time: 2020-07-07 17:44:18 -0400



Basin Characteristics					
Parameter Code	Parameter Description	Value	Unit		
STRMTOT	total length of all mapped streams (1:24,000-scale) in the basin	6.26	miles		
DRNAREA	Area that drains to a point on a stream	3.63	square miles		
BSLDEM250	Mean basin slope computed from 1:250K DEM	1.483	percent		
DRFTPERSTR	Area of stratified drift per unit of stream length	0.3	square mile per mile		
MAREGION	Region of Massachusetts 0 for Eastern 1 for Western	0	dimensionless		

https://streamstats.usgs.gov/ss/

Parameter Code	Parameter Description	Value	Unit
PCTSNDGRV	Percentage of land surface underlain by sand and gravel deposits	51.62	percent
FOREST	Percentage of area covered by forest	16.95	percent
BSLDEM10M	Mean basin slope computed from 10 m DEM	5.01	percent
ELEV	Mean Basin Elevation	218	feet
LC06STOR	Percentage of water bodies and wetlands determined from the NLCD 2006	5.25	percent

Low-Flow Statistics Parameters [Statewide Low Flow WRIR00 4135]

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63	square miles	1.61	149
BSLDEM250	Mean Basin Slope from 250K DEM	1.483	percent	0.32	24.6
DRFTPERSTR	Stratified Drift per Stream Length	0.3	square mile per mile	0	1.29
MAREGION	Massachusetts Region	0	dimensionless	0	1

Low-Flow Statistics Flow Report[Statewide Low Flow WRIR00 4135]

PII: Prediction Interval-Lower, Plu: Prediction Interval-Upper, SEp: Standard Error of Prediction, SE: Standard Error (other -- see report)

Statistic	Value	Unit	PII	Plu	SE	SEp
7 Day 2 Year Low Flow	0.377	ft^3/s	0.13	1.05	49.5	49.5
7 Day 10 Year Low Flow	0.15	ft^3/s	0.0405	0.518	70.8	70.8

Low-Flow Statistics Citations

Ries, K.G., III,2000, Methods for estimating low-flow statistics for Massachusetts streams: U.S. Geological Survey Water Resources Investigations Report 00-4135, 81 p. (http://pubs.usgs.gov/wri/wri004135/)

Flow-Duration Statistics Parameters[Statewide Low Flow WRIR00 4135]

https://streamstats.usgs.gov/ss/ 2/7

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63	square miles	1.61	149
DRFTPERSTR	Stratified Drift per Stream Length	0.3	square mile per mile	0	1.29
MAREGION	Massachusetts Region	0	dimensionless	0	1
BSLDEM250	Mean Basin Slope from 250K DEM	1.483	percent	0.32	24.6

Flow-Duration Statistics Flow Report[Statewide Low Flow WRIR00 4135]

PII: Prediction Interval-Lower, Plu: Prediction Interval-Upper, SEp: Standard Error of Prediction, SE: Standard Error (other -- see report)

Statistic	Value	Unit	PII	Plu	SE	SEp
50 Percent Duration	3.56	ft^3/s	1.62	7.74	17.6	17.6
60 Percent Duration	2.64	ft^3/s	1.02	6.77	19.8	19.8
70 Percent Duration	1.74	ft^3/s	0.748	4	23.5	23.5
75 Percent Duration	1.38	ft^3/s	0.6	3.13	25.8	25.8
80 Percent Duration	1.14	ft^3/s	0.509	2.5	28.4	28.4
85 Percent Duration	0.82	ft^3/s	0.343	1.92	31.9	31.9
90 Percent Duration	0.628	ft^3/s	0.256	1.5	36.6	36.6
95 Percent Duration	0.347	ft^3/s	0.123	0.951	45.6	45.6
98 Percent Duration	0.228	ft^3/s	0.0704	0.698	60.3	60.3
99 Percent Duration	0.165	ft^3/s	0.0478	0.538	65.1	65.1

Flow-Duration Statistics Citations

Ries, K.G., III,2000, Methods for estimating low-flow statistics for Massachusetts streams: U.S. Geological Survey Water Resources Investigations Report 00-4135, 81 p. (http://pubs.usgs.gov/wri/wri004135/)

August Flow-Duration Statistics Parameters[Statewide Low Flow WRIR00 4135]

Parameter Code	Parameter Name	Value Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63 square miles	1.61	149

https://streamstats.usgs.gov/ss/ 3/7

StreamStats

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
BSLDEM250	Mean Basin Slope from 250K DEM	1.483	percent	0.32	24.6
DRFTPERSTR	Stratified Drift per Stream Length	0.3	square mile per mile	0	1.29
MAREGION	Massachusetts Region	0	dimensionless	0	1

August Flow-Duration Statistics Flow Report[Statewide Low Flow WRIR00 4135]

PII: Prediction Interval-Lower, Plu: Prediction Interval-Upper, SEp: Standard Error of Prediction, SE: Standard Error (other -- see report)

Statistic	Value	Unit	PII	Plu	SE	SEp
August 50 Percent Duration	0.911	ft^3/s	0.37	2.2	33.2	33.2

August Flow-Duration Statistics Citations

Ries, K.G., III,2000, Methods for estimating low-flow statistics for Massachusetts streams: U.S. Geological Survey Water Resources Investigations Report 00-4135, 81 p. (http://pubs.usgs.gov/wri/wri004135/)

Probability Statistics Parameters[Perennial Flow Probability]

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63	square miles	0.01	1.99
PCTSNDGRV	Percent Underlain By Sand And Gravel	51.62	percent	0	100
FOREST	Percent Forest	16.95	percent	0	100
MAREGION	Massachusetts Region	0	dimensionless	0	1

Probability Statistics Disclaimers[Perennial Flow Probability]

One or more of the parameters is outside the suggested range. Estimates were extrapolated with unknown errors

Probability Statistics Flow Report[Perennial Flow Probability]

Statistic	Value	Unit
Probability Stream Flowing Perennially	0.986	dim

Probability Statistics Citations

Bent, G.C., and Steeves, P.A.,2006, A revised logistic regression equation and an automated procedure for mapping the probability of a stream flowing perennially in Massachusetts: U.S. Geological Survey Scientific Investigations Report 2006–5031, 107 p. (http://pubs.usgs.gov/sir/2006/5031/pdfs/SIR_2006-5031rev.pdf)

Bankfull Statistics Parameters[Bankfull Statewide SIR2013 5155]

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63	square miles	0.6	329
BSLDEM10M	Mean Basin Slope from 10m DEM	5.01	percent	2.2	23.9

Bankfull Statistics Flow Report[Bankfull Statewide SIR2013 5155]

PII: Prediction Interval-Lower, Plu: Prediction Interval-Upper, SEp: Standard Error of Prediction, SE: Standard Error (other -- see report)

Statistic	Value	Unit	SEp
Bankfull Width	23.5	ft	21.3
Bankfull Depth	1.31	ft	19.8
Bankfull Area	30.6	ft^2	29
Bankfull Streamflow	75	ft^3/s	55

Bankfull Statistics Citations

Bent, G.C., and Waite, A.M.,2013, Equations for estimating bankfull channel geometry and discharge for streams in Massachusetts: U.S. Geological Survey Scientific Investigations Report 2013–5155, 62 p., (http://pubs.usgs.gov/sir/2013/5155/)

Peak-Flow Statistics Parameters[Peak Statewide 2016 5156]

Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit
DRNAREA	Drainage Area	3.63	square miles	0.16	512
ELEV	Mean Basin Elevation	218	feet	80.6	1948

https://streamstats.usgs.gov/ss/ 5/7

Parameter Code	Parameter Name	Value Units	Min Limit	Max Limit
LC06STOR	Percent Storage from NLCD2006	5.25 percent	0	32.3

Peak-Flow Statistics Flow Report[Peak Statewide 2016 5156]

PII: Prediction Interval-Lower, Plu: Prediction Interval-Upper, SEp: Standard Error of Prediction, SE: Standard Error (other -- see report)

Statistic	Value	Unit	PII	Plu	SEp
2 Year Peak Flood	115	ft^3/s	58.4	224	42.3
5 Year Peak Flood	189	ft^3/s	95	375	43.4
10 Year Peak Flood	247	ft^3/s	121	503	44.7
25 Year Peak Flood	332	ft^3/s	158	699	47.1
50 Year Peak Flood	402	ft^3/s	185	875	49.4
100 Year Peak Flood	477	ft^3/s	212	1070	51.8
200 Year Peak Flood	558	ft^3/s	241	1290	54.1
500 Year Peak Flood	675	ft^3/s	279	1630	57.6

Peak-Flow Statistics Citations

Zarriello, P.J.,2017, Magnitude of flood flows at selected annual exceedance probabilities for streams in Massachusetts: U.S. Geological Survey Scientific Investigations Report 2016-5156, 99 p. (https://dx.doi.org/10.3133/sir20165156)

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Application Version: 4.3.11

ATTACHMENT C MACRIS REPORT

Massachusetts Cultural Resource Information System MACRIS

MACRIS Search Results

Search Criteria: Town(s): Burlington; Street Name: Wheeler; Resource Type(s): Area, Building, Burial Ground, Object, Structure;

Inv. No. Property Name Street Town Year

Thursday, June 4, 2020 Page 1 of 1

ATTACHMENT D ENDANGERED AND CRITICAL SPECIES REPORT

IPaC

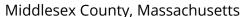
U.S. Fish & Wildlife Service

IPaC resource list

This report is an automatically generated list of species and other resources such as critical habitat (collectively referred to as *trust resources*) under the U.S. Fish and Wildlife Service's (USFWS) jurisdiction that are known or expected to be on or near the project area referenced below. The list may also include trust resources that occur outside of the project area, but that could potentially be directly or indirectly affected by activities in the project area. However, determining the likelihood and extent of effects a project may have on trust resources typically requires gathering additional site-specific (e.g., vegetation/species surveys) and project-specific (e.g., magnitude and timing of proposed activities) information.

Below is a summary of the project information you provided and contact information for the USFWS office(s) with jurisdiction in the defined project area. Please read the introduction to each section that follows (Endangered Species, Migratory Birds, USFWS Facilities, and NWI Wetlands) for additional information applicable to the trust resources addressed in that section.

Location





Local office

New England Ecological Services Field Office

(603) 223-2541

(603) 223-0104

70 Commercial Street, Suite 300 Concord, NH 03301-5094

http://www.fws.gov/newengland

Endangered species

This resource list is for informational purposes only and does not constitute an analysis of project level impacts.

The primary information used to generate this list is the known or expected range of each species. Additional areas of influence (AOI) for species are also considered. An AOI includes areas outside of the species range if the species could be indirectly affected by activities in that area (e.g., placing a dam upstream of a fish population, even if that fish does not occur at the dam site, may indirectly impact the species by reducing or eliminating water flow downstream). Because species can move, and site conditions can change, the species on this list are not guaranteed to be found on or near the project area. To fully determine any potential effects to species, additional site-specific and project-specific information is often required.

Section 7 of the Endangered Species Act **requires** Federal agencies to "request of the Secretary information whether any species which is listed or proposed to be listed may be present in the area of such proposed action" for any project that is conducted, permitted, funded, or licensed by any Federal agency. A letter from the local office and a species list which fulfills this requirement can **only** be obtained by requesting an official species list from either the Regulatory Review section in IPaC (see directions below) or from the local field office directly.

For project evaluations that require USFWS concurrence/review, please return to the IPaC website and request an official species list by doing the following:

- 1. Draw the project location and click CONTINUE.
- 2. Click DEFINE PROJECT.
- 3. Log in (if directed to do so).
- 4. Provide a name and description for your project.
- 5. Click REQUEST SPECIES LIST.

Listed species¹ and their critical habitats are managed by the <u>Ecological Services Program</u> of the U.S. Fish and Wildlife Service (USFWS) and the fisheries division of the National Oceanic and Atmospheric Administration (NOAA Fisheries²).

Species and critical habitats under the sole responsibility of NOAA Fisheries are **not** shown on this list. Please contact <u>NOAA Fisheries</u> for <u>species under their jurisdiction</u>.

- 1. Species listed under the <u>Endangered Species Act</u> are threatened or endangered; IPaC also shows species that are candidates, or proposed, for listing. See the <u>listing status page</u> for more information.
- 2. <u>NOAA Fisheries</u>, also known as the National Marine Fisheries Service (NMFS), is an office of the National Oceanic and Atmospheric Administration within the Department of Commerce.

The following species are potentially affected by activities in this location:

Mammals

NAME STATUS

Northern Long-eared Bat Myotis septentrionalis No critical habitat has been designated for this species. https://ecos.fws.gov/ecp/species/9045 **Threatened**

Critical habitats

Potential effects to critical habitat(s) in this location must be analyzed along with the endangered species themselves.

THERE ARE NO CRITICAL HABITATS AT THIS LOCATION.

Migratory birds

Certain birds are protected under the Migratory Bird Treaty Act^{1} and the Bald and Golden Eagle Protection Act^{2} .

Any person or organization who plans or conducts activities that may result in impacts to migratory birds, eagles, and their habitats should follow appropriate regulations and consider implementing appropriate conservation measures, as described <u>below</u>.

- 1. The Migratory Birds Treaty Act of 1918.
- 2. The Bald and Golden Eagle Protection Act of 1940.

Additional information can be found using the following links:

- Birds of Conservation Concern http://www.fws.gov/birds/management/managed-species/birds-of-conservation-concern.php
- Measures for avoiding and minimizing impacts to birds
 http://www.fws.gov/birds/management/project-assessment-tools-and-guidance/conservation-measures.php
- Nationwide conservation measures for birds http://www.fws.gov/migratorybirds/pdf/management/nationwidestandardconservationmeasures.pdf

The birds listed below are birds of particular concern either because they occur on the <u>USFWS Birds of Conservation Concern</u> (BCC) list or warrant special attention in your project location. To learn more about the levels of concern for birds on your list and how this list is generated, see the FAQ <u>below</u>. This is not a list of every bird you may find in this location, nor a guarantee that every bird on this list will be found in your project area. To see exact locations of where birders and the general public have sighted birds in and around your project area, visit the <u>E-bird data mapping tool</u> (Tip: enter your location, desired date range and a species on your list). For projects that occur off the Atlantic Coast, additional maps and models detailing the relative occurrence and abundance of bird species on your list are available. Links to additional information about Atlantic Coast birds, and other important information about your migratory bird list, including how to properly interpret and use your migratory bird report, can be found <u>below</u>.

IPaC: Explore Location

For guidance on when to schedule activities or implement avoidance and minimization measures to reduce impacts to migratory birds on your list, click on the PROBABILITY OF PRESENCE SUMMARY at the top of your list to see when these birds are most likely to be present and breeding in your project area.

NAME

BREEDING SEASON (IF A
BREEDING SEASON IS INDICATED
FOR A BIRD ON YOUR LIST, THE
BIRD MAY BREED IN YOUR
PROJECT AREA SOMETIME WITHIN
THE TIMEFRAME SPECIFIED,
WHICH IS A VERY LIBERAL
ESTIMATE OF THE DATES INSIDE
WHICH THE BIRD BREEDS
ACROSS ITS ENTIRE RANGE.
"BREEDS ELSEWHERE" INDICATES
THAT THE BIRD DOES NOT LIKELY
BREED IN YOUR PROJECT AREA.)

Bald Eagle Haliaeetus leucocephalus

This is not a Bird of Conservation Concern (BCC) in this area, but warrants attention because of the Eagle Act or for potential susceptibilities in offshore areas from certain types of development or activities.

https://ecos.fws.gov/ecp/species/1626

Black-billed Cuckoo Coccyzus erythropthalmus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

https://ecos.fws.gov/ecp/species/9399

Bobolink Dolichonyx oryzivorus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Buff-breasted Sandpiper Calidris subruficollis

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

https://ecos.fws.gov/ecp/species/9488

Canada Warbler Cardellina canadensis

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Dunlin Calidris alpina arcticola

This is a Bird of Conservation Concern (BCC) only in particular Bird Conservation Regions (BCRs) in the continental USA

Breeds Oct 15 to Aug 31

Breeds May 15 to Oct 10

Breeds May 20 to Jul 31

Breeds elsewhere

Breeds May 20 to Aug 10

Breeds elsewhere

IPaC: Explore Location

Evening Grosbeak Coccothraustes vespertinus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds elsewhere

Lesser Yellowlegs Tringa flavipes

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

https://ecos.fws.gov/ecp/species/9679

Breeds elsewhere

Prairie Warbler Dendroica discolor

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds May 1 to Jul 31

Red-headed Woodpecker Melanerpes erythrocephalus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds May 10 to Sep 10

Rusty Blackbird Euphagus carolinus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds elsewhere

Semipalmated Sandpiper Calidris pusilla

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds elsewhere

Short-billed Dowitcher Limnodromus griseus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

https://ecos.fws.gov/ecp/species/9480

Breeds elsewhere

Snowy Owl Bubo scandiacus

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds elsewhere

Wood Thrush Hylocichla mustelina

This is a Bird of Conservation Concern (BCC) throughout its range in the continental USA and Alaska.

Breeds May 10 to Aug 31

Probability of Presence Summary

The graphs below provide our best understanding of when birds of concern are most likely to be present in your project area. This information can be used to tailor and schedule your project activities to avoid or minimize impacts to birds. Please make sure you read and understand the FAQ "Proper Interpretation and Use of Your Migratory Bird Report" before using or attempting to interpret this report.

Probability of Presence (■)

Each green bar represents the bird's relative probability of presence in the 10km grid cell(s) your project overlaps during a particular week of the year. (A year is represented as 12 4-week months.) A taller bar indicates a higher probability of species presence. The survey effort (see below) can be used to establish a level of confidence in the presence score. One can have higher confidence in the presence score if the corresponding survey effort is also high.

How is the probability of presence score calculated? The calculation is done in three steps:

- 1. The probability of presence for each week is calculated as the number of survey events in the week where the species was detected divided by the total number of survey events for that week. For example, if in week 12 there were 20 survey events and the Spotted Towhee was found in 5 of them, the probability of presence of the Spotted Towhee in week 12 is 0.25.
- 2. To properly present the pattern of presence across the year, the relative probability of presence is calculated. This is the probability of presence divided by the maximum probability of presence across all weeks. For example, imagine the probability of presence in week 20 for the Spotted Towhee is 0.05, and that the probability of presence at week 12 (0.25) is the maximum of any week of the year. The relative probability of presence on week 12 is 0.25/0.25 = 1; at week 20 it is 0.05/0.25 = 0.2.
- 3. The relative probability of presence calculated in the previous step undergoes a statistical conversion so that all possible values fall between 0 and 10, inclusive. This is the probability of presence score.

To see a bar's probability of presence score, simply hover your mouse cursor over the bar.

Breeding Season (

Yellow bars denote a very liberal estimate of the time-frame inside which the bird breeds across its entire range. If there are no yellow bars shown for a bird, it does not breed in your project area.

Survey Effort (1)

Vertical black lines superimposed on probability of presence bars indicate the number of surveys performed for that species in the 10km grid cell(s) your project area overlaps. The number of surveys is expressed as a range, for example, 33 to 64 surveys.

To see a bar's survey effort range, simply hover your mouse cursor over the bar.

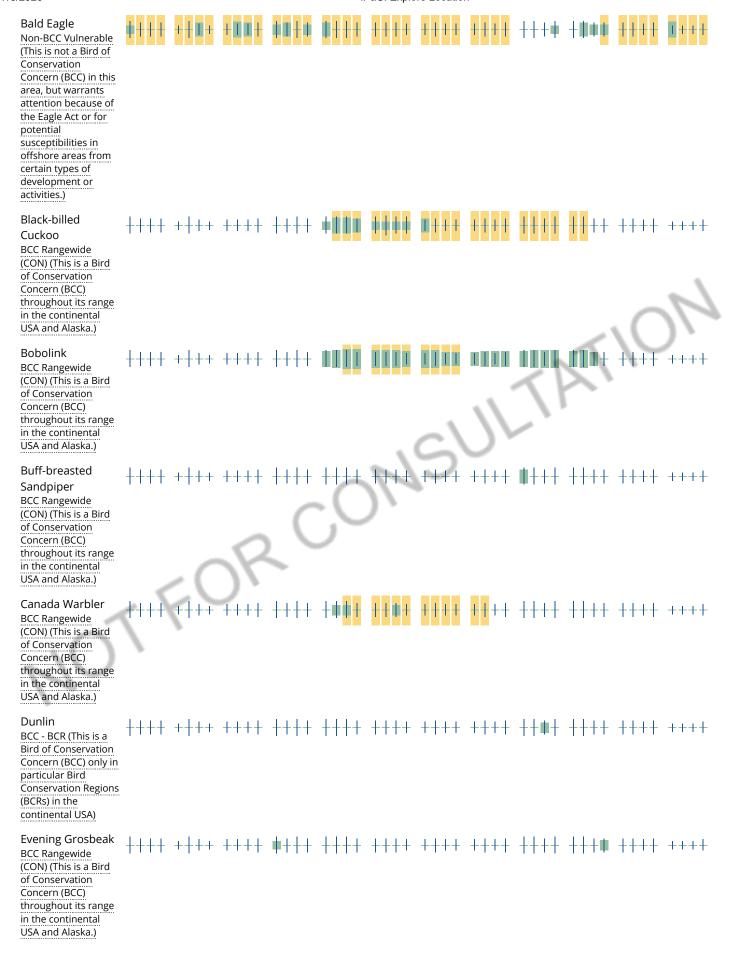
No Data (-)

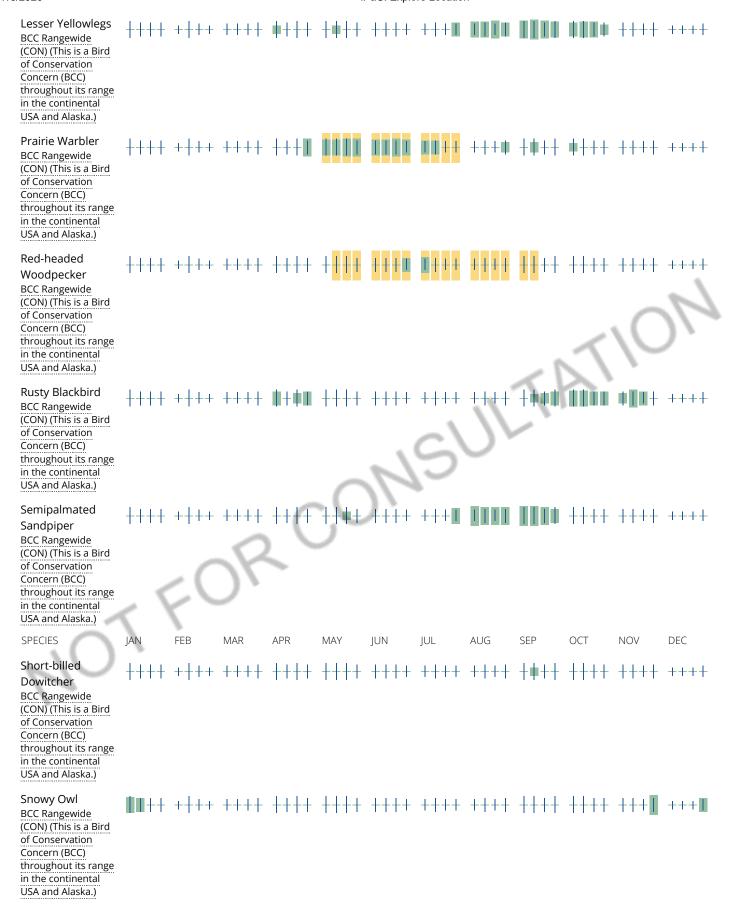
A week is marked as having no data if there were no survey events for that week.

Survey Timeframe

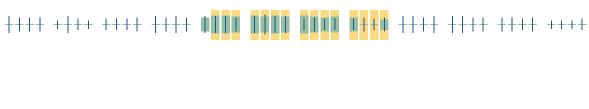
Surveys from only the last 10 years are used in order to ensure delivery of currently relevant information. The exception to this is areas off the Atlantic coast, where bird returns are based on all years of available data, since data in these areas is currently much more sparse.











Tell me more about conservation measures I can implement to avoid or minimize impacts to migratory birds.

Nationwide Conservation Measures describes measures that can help avoid and minimize impacts to all birds at any location year round. Implementation of these measures is particularly important when birds are most likely to occur in the project area. When birds may be breeding in the area, identifying the locations of any active nests and avoiding their destruction is a very helpful impact minimization measure. To see when birds are most likely to occur and be breeding in your project area, view the Probability of Presence Summary. Additional measures and/or permits may be advisable depending on the type of activity you are conducting and the type of infrastructure or bird species present on your project site.

What does IPaC use to generate the migratory birds potentially occurring in my specified location?

The Migratory Bird Resource List is comprised of USFWS <u>Birds of Conservation Concern (BCC)</u> and other species that may warrant special attention in your project location.

The migratory bird list generated for your project is derived from data provided by the <u>Avian Knowledge Network (AKN)</u>. The AKN data is based on a growing collection of <u>survey, banding, and citizen science datasets</u> and is queried and filtered to return a list of those birds reported as occurring in the 10km grid cell(s) which your project intersects, and that have been identified as warranting special attention because they are a BCC species in that area, an eagle (<u>Eagle Act</u> requirements may apply), or a species that has a particular vulnerability to offshore activities or development.

Again, the Migratory Bird Resource list includes only a subset of birds that may occur in your project area. It is not representative of all birds that may occur in your project area. To get a list of all birds potentially present in your project area, please visit the AKN Phenology Tool.

What does IPaC use to generate the probability of presence graphs for the migratory birds potentially occurring in my specified location?

The probability of presence graphs associated with your migratory bird list are based on data provided by the <u>Avian Knowledge Network (AKN)</u>. This data is derived from a growing collection of <u>survey</u>, <u>banding</u>, <u>and citizen</u> science datasets .

Probability of presence data is continuously being updated as new and better information becomes available. To learn more about how the probability of presence graphs are produced and how to interpret them, go the Probability of Presence Summary and then click on the "Tell me about these graphs" link.

How do I know if a bird is breeding, wintering, migrating or present year-round in my project area?

To see what part of a particular bird's range your project area falls within (i.e. breeding, wintering, migrating or year-round), you may refer to the following resources: The Cornell Lab of Ornithology All About Birds Bird Guide, or (if you are unsuccessful in locating the bird of interest there), the Cornell Lab of Ornithology Neotropical Birds guide. If a bird on your migratory bird species list has a breeding season associated with it, if that bird does occur in your project area, there may be nests present at some point within the timeframe specified. If "Breeds elsewhere" is indicated, then the bird likely does not breed in your project area.

What are the levels of concern for migratory birds?

Migratory birds delivered through IPaC fall into the following distinct categories of concern:

- 1. "BCC Rangewide" birds are <u>Birds of Conservation Concern</u> (BCC) that are of concern throughout their range anywhere within the USA (including Hawaii, the Pacific Islands, Puerto Rico, and the Virgin Islands);
- 2. "BCC BCR" birds are BCCs that are of concern only in particular Bird Conservation Regions (BCRs) in the continental USA; and
- 3. "Non-BCC Vulnerable" birds are not BCC species in your project area, but appear on your list either because of the <u>Eagle Act</u> requirements (for eagles) or (for non-eagles) potential susceptibilities in offshore areas from certain types of development or activities (e.g. offshore energy development or longline fishing).

Although it is important to try to avoid and minimize impacts to all birds, efforts should be made, in particular, to avoid and minimize impacts to the birds on this list, especially eagles and BCC species of rangewide concern. For more information on conservation measures you can implement to help avoid and minimize migratory bird impacts and requirements for eagles, please see the FAQs for these topics.

Details about birds that are potentially affected by offshore projects

For additional details about the relative occurrence and abundance of both individual bird species and groups of bird species within your project area off the Atlantic Coast, please visit the <u>Northeast Ocean Data Portal</u>. The Portal also offers data and information about other taxa besides birds that may be helpful to you in your project review. Alternately, you may download the bird model results files underlying the portal maps through the <u>NOAA NCCOS</u> <u>Integrative Statistical Modeling and Predictive Mapping of Marine Bird Distributions and Abundance on the Atlantic Outer Continental Shelf</u> project webpage.

Bird tracking data can also provide additional details about occurrence and habitat use throughout the year, including migration. Models relying on survey data may not include this information. For additional information on marine bird tracking data, see the <u>Diving Bird Study</u> and the <u>nanotag studies</u> or contact <u>Caleb Spiegel</u> or <u>Pam Loring</u>.

What if I have eagles on my list?

If your project has the potential to disturb or kill eagles, you may need to <u>obtain a permit</u> to avoid violating the Eagle Act should such impacts occur.

Proper Interpretation and Use of Your Migratory Bird Report

The migratory bird list generated is not a list of all birds in your project area, only a subset of birds of priority concern. To learn more about how your list is generated, and see options for identifying what other birds may be in your project area, please see the FAQ "What does IPaC use to generate the migratory birds potentially occurring in my specified location". Please be aware this report provides the "probability of presence" of birds within the 10 km grid cell(s) that overlap your project; not your exact project footprint. On the graphs provided, please also look carefully at the survey effort (indicated by the black vertical bar) and for the existence of the "no data" indicator (a red horizontal bar). A high survey effort is the key component. If the survey effort is high, then the probability of presence score can be viewed as more dependable. In contrast, a low survey effort bar or no data bar means a lack of data and, therefore, a lack of certainty about presence of the species. This list is not perfect; it is simply a starting point for identifying what birds of concern have the potential to be in your project area, when they might be there, and if they might be breeding (which means nests might be present). The list helps you know what to look for to confirm presence, and helps guide you in knowing when to implement conservation measures to avoid or minimize potential impacts from your project activities, should presence be confirmed. To learn more about conservation measures, visit the FAQ "Tell me about conservation measures I can implement to avoid or minimize impacts to migratory birds" at the bottom of your migratory bird trust resources page.



National Wildlife Refuge lands

Any activity proposed on lands managed by the <u>National Wildlife Refuge</u> system must undergo a 'Compatibility Determination' conducted by the Refuge. Please contact the individual Refuges to discuss any questions or concerns.

THERE ARE NO REFUGE LANDS AT THIS LOCATION.

Fish hatcheries

THERE ARE NO FISH HATCHERIES AT THIS LOCATION.

Wetlands in the National Wetlands Inventory

Impacts to <u>NWI wetlands</u> and other aquatic habitats may be subject to regulation under Section 404 of the Clean Water Act, or other State/Federal statutes.

For more information please contact the Regulatory Program of the local <u>U.S. Army Corps of Engineers District</u>.

Please note that the NWI data being shown may be out of date. We are currently working to update our NWI data set. We recommend you verify these results with a site visit to determine the actual extent of wetlands on site.

This location overlaps the following wetlands:

FRESHWATER EMERGENT WETLAND

PEM1E

PEM1Fh

RIVERINE

R2UBH

R2UBHx

A full description for each wetland code can be found at the National Wetlands Inventory website

Data limitations

The Service's objective of mapping wetlands and deepwater habitats is to produce reconnaissance level information on the location, type and size of these resources. The maps are prepared from the analysis of high altitude imagery. Wetlands are identified based on vegetation, visible hydrology and geography. A margin of error is inherent in the use of imagery; thus, detailed on-the-ground inspection of any particular site may result in revision of the wetland boundaries or classification established through image analysis.

The accuracy of image interpretation depends on the quality of the imagery, the experience of the image analysts, the amount and quality of the collateral data and the amount of ground truth verification work conducted. Metadata should be consulted to determine the date of the source imagery used and any mapping problems.

Wetlands or other mapped features may have changed since the date of the imagery or field work. There may be occasional differences in polygon boundaries or classifications between the information depicted on the map and the actual conditions on site.

Data exclusions

Certain wetland habitats are excluded from the National mapping program because of the limitations of aerial imagery as the primary data source used to detect wetlands. These habitats include seagrasses or submerged aquatic vegetation that are found in the intertidal and subtidal zones of estuaries and nearshore coastal waters. Some deepwater reef communities (coral or tuberficid worm reefs) have also been excluded from the inventory. These habitats, because of their depth, go undetected by aerial imagery.

Data precautions

Federal, state, and local regulatory agencies with jurisdiction over wetlands may define and describe wetlands in a different manner than that used in this inventory. There is no attempt, in either the design or products of this inventory, to define the limits of proprietary jurisdiction of any Federal, state, or local government or to establish the geographical scope of the regulatory programs of government agencies. Persons intending to engage in activities involving modifications within or adjacent to wetland areas should seek the advice of appropriate federal, state, or local agencies concerning specified agency regulatory programs and proprietary jurisdictions that may affect such activities.

ATTACHMENT E EFFLUENT DATA ESTIMATIONS

PROJECT EFFLUENT CONCENTRATIONS

Water Treatment System Wheeler Road Burlington, MA

Influent Contaminant	Highest Influent Concentration (ug/L)	Treatment Method	Removal Efficiency ² (%)	Estimated Effluent Concentration (ug/L)
1,2,4-Trimethylbenzene	8.4	Carbon	99	0.08
1,3,5-Trimethylbenzene	3.3	Carbon	99	0.03
Dichlorodifluoromethane	7.1	Carbon	99	0.07
o-Xylene	2.9	Carbon	99	0.03
m&p-Xylene	4.3	Carbon	99	0.04
Naphthalene	2.7	Carbon	99	0.03
Vinyl Chloride	2	Carbon	99 ³	0.02
cis 1,2-dichloroethane	30	Carbon	99	0.3
Trichloroethene (TCE)	62	Carbon	95	3.1
Tetrachloroethene (PCE)	18	Carbon	95	0.9
Antimony	10	Settling, Filtration, Ion Exchange	45	5.5
Arsenic	7	Settling, Filtration, Ion Exchange	80	1.4
Barium	106	Settling, Filtration, Ion Exchange	85	15.9
Chromium	12	Settling, Filtration, Ion Exchange	97	0.36
Lead	9.0	Settling, Filtration, Ion Exchange	99	0.09
Nickel	8.0	Settling, Filtration, Ion Exchange	84	1.28
Vanadium	9.0	Settling, Filtration, Ion Exchange	85	1.35
Zinc	30	Settling, Filtration, Ion Exchange	82	5.4

Notes

- 1. Treatment system includes an 18,000-gallon weir tank, centrifugal transfer pump, duplex bag filter skid, two 6,000 lb carbon vessels in series, a 120 CF cation exchange vessel, and a flow/totalizer meter
- 2. Removal efficiency based on empirical data
- 3. Removal efficiency of vinyl chloride may readily decrease depending on loading of other CVOC and VOC contaminants

ATTACHMENT F GROUNDWATER DATA

TABLE 2

Ground Water and Surface Water Analytical Data March 2017 and March 2020 Wheeler Rd., Burlington, Massachusetts BL Companies Project No. 16D3253

Compounds	MCP Reportable Concentration for Groundwater		Concentration of Compound in Sample												
•	RCGW-1	MW-1	MW-1/DUP	MW-1	MW-2	MW-2	MW-3	MW-4	MW-101	MW-102	MW-103	MW-104	SW-1	SW-2	Trip Blank
		3/27/2017	3/27/2017	3/17/2020	3/27/2017	3/17/2020	3/27/2017	3/27/2017	3/17/2020	3/17/2020	3/17/2020	3/17/2020	3/27/2017	3/27/2014	3/27/2017
MADEP-VPH (ug/L)															
Unadjusted C5-C8 Aliphatics	300	ND	ND	NA	ND	ND <100	ND	ND	ND <100	NA	ND <100	NA	ND	ND	ND
C5-C8 Aliphatic Hydrocarbons	300	ND	ND	NA	ND	ND <100	ND	ND	ND <100	NA	ND <100	NA	ND	ND	ND
Unadjusted C9-C12 Aliphatics	700	ND	ND	NA	ND	ND <100	ND	ND	150	NA	ND <100	NA	ND	ND	ND
C9-C12 Aliphatic Hydrocarbons	700	ND	ND	NA	ND	ND <100	ND	ND	ND <100	NA	ND <100	NA	ND	ND	ND
C9-C10 Aromatic Hydrocarbons	200	ND	ND	NA	ND	ND <100	ND	ND	ND <100	NA	ND <100	NA	ND	ND	ND
MADEP-EPH (μg/L)															
Unadusted C11-C22 Aromatic Hydrocarbons	200	ND	ND	NA	ND	ND <200	ND	ND	ND <200	NA	ND <200	NA	ND	ND	NA
C11-C22 Aromatic Hydrocarbons	200	ND	ND	NA	ND	ND <200	ND	ND	ND <200	NA	ND <200	NA	ND	ND	NA
C9-C18 Alphatic Hydrocarbons	700	ND	ND	NA	ND	ND <200	ND	ND	ND <200	NA	ND <200	NA	ND	ND	NA
C19-C36 Aliphatic Hydrocarbons	14,000	ND	ND	NA	ND	ND <200	ND	ND	ND <200	NA	ND <200	NA	ND	ND	NA
SVOCs (μg/L)	Varies	NA	NA	NA	NA	ND	NA	NA	ND	NA	ND	NA	NA	NA	NA
/OCs (μg/L)															
1.2.4-Trimethylbenzene	10.000	ND	ND	ND	ND	ND	ND	ND	8.1	ND	ND	ND	ND	ND	ND
1.3.5-Trimethylbenzene	100	ND	ND	ND	ND	ND	ND	ND	3.3	ND	ND	ND	ND	ND	ND
Dichlorodifluoromethane	10,000	ND	ND	ND	ND	ND	ND	ND	ND	2.6	7.1	ND	ND	ND	ND
o-Xvlene	6,000	ND	ND	ND	ND	ND	ND	ND	2.9	ND.	ND	ND	ND	ND	ND
m&p-Xylene	4,000	ND	ND	ND	ND	ND	ND	ND	4.3	ND	ND	ND	ND	ND	ND
Total Xvlene	500	ND	ND	ND	ND	ND	ND	ND	7.2	ND	ND	ND	ND	ND	ND
Naphthalene	140	ND	ND	ND	ND	ND	ND	ND	2.7	ND	ND	ND	ND	ND	ND
Vinyl Chloride	2	2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1.1-Dichloroethene	7	ND	ND	ND	2	ND									
cis 1.2-dichloroethane	20	2	3	1.7	14	3.7	ND	ND	6.4	5.4	30	ND	ND	ND	ND
Trichloroethene (TCE)	5	5	4	1.7	62	20	ND	ND	44	5.8	34	ND	2	ND	ND
Tetrachloroethene (PCE)	5	1	ND	ND	18	4.5	ND	ND	7	ND	5.4	ND	ND	ND	ND
Total MCP 14 Metals (μg/L)															+
Antimony	6	10	ND	NA	9	ND < 5	10	7	ND < 5	NA	ND < 5	NA	ND	ND	NA
Arsenic	10	ND	ND	NA	ND	ND < 4	ND	ND	7	NA	ND <4	NA	ND	ND	NA
Barium	100	95	89	NA	94	106	ND	40	99	NA	13	NA	49	31	NA
Chromium	100	ND	ND	NA	ND	2	ND	ND	12	NA	ND <1	NA	ND	ND	NA
_ead	10	ND	ND	NA	ND	ND <2	ND	ND	9	NA	ND < 2	NA	ND	6	NA
Nickel	100	ND	ND	NA	ND	4	ND	ND	8	NA	ND <1	NA	ND	ND	NA
/anadium	30	ND	ND	NA	ND	3	ND	ND	9	NA	ND <2	NA	ND	ND	NA
Zinc	900	20	ND	NA	30	17	ND	ND	17	NA	ND <4	NA	30	ND	NA
Dissolved MCP 14 Metals (µg/L)															<u> </u>
Antimony	6	ND	NA	NA	ND	NA	ND	ND	NA						

NOTES:

NOTES:
Only compounds detected are listed
Dissolved antimony samples were collected on May 18, 2017

µg/L = micrograms per liter = parts per billion

ND = Not detected

ND = Not detected

NA= Not anaylzed Bold and shaded values indicate an exceedence of the MCP RCGW-1



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

Analysis Report

May 14, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: GROUND WATER Collected by: 05/05/20

Location Code: BLCOMPANIES Received by: B 05/06/20 15:46

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF90108

Phoenix ID: CF90196

Project ID: 16D3253 Client ID: MW-1 25UM

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	05/07/20	TH	SW6010D
Barium	0.058	0.002	mg/L	1	05/07/20	TH	SW6010D
Beryllium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Chromium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Mercury	< 0.0002	0.0002	mg/L	1	05/08/20	RS	SW7470A
Nickel	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Lead	< 0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Antimony	< 0.005	0.005	mg/L	1	05/07/20	TH	SW6010D
Selenium	< 0.010	0.010	mg/L	1	05/07/20	TH	SW6010D
Thallium	< 0.001	0.001	mg/L	1	05/07/20	RS	SW7010
Vanadium	< 0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Zinc	< 0.004	0.004	mg/L	1	05/07/20	TH	SW6010D
Filtration	Completed					/BF	0.45um Filter
Mercury Digestion	Completed				05/07/20	RA/RA	SW7470A
Total Metals Digestion	Completed				05/06/20	AG/BF	•

Client ID: MW-1 25UM

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

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

Phyllis Shiller, Laboratory Director

May 14, 2020

Reviewed and Released by: Phyllis Shiller, Laboratory Director



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

Analysis Report

May 14, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Phoenix ID: CF90197

<u>Sample Information</u> <u>Custody Information</u> <u>Date</u> <u>Time</u>

Matrix: GROUND WATER Collected by: 05/05/20

Location Code: BLCOMPANIES Received by: B 05/06/20 15:46

Rush Request: 72 Hour Analyzed by: see "By" below

Laboratory Data SDG ID: GCF90108

Project ID: 16D3253 Client ID: MW-2 25UM

P.O.#:

RL/ Parameter Result **PQL** Units Dilution Date/Time By Reference Silver < 0.001 0.001 mg/L 1 05/07/20 TH SW6010D Arsenic < 0.004 0.004 mg/L 1 05/07/20 TH SW6010D 05/07/20 SW6010D Barium 0.094 0.002 mg/L 1 TH < 0.001 Beryllium 0.001 mg/L 1 05/07/20 ΤH SW6010D SW6010D < 0.001 0.001 mg/L 1 05/07/20 TH Cadmium Chromium 0.009 0.001 mg/L 1 05/07/20 TH SW6010D < 0.0002 0.0002 05/08/20 SW7470A Mercury mg/L 1 RS 0.012 0.001 mg/L 1 05/07/20 TH SW6010D Nickel 05/07/20 SW6010D 0.005 0.002 mg/L 1 ΤH Lead SW6010D **Antimony** < 0.005 0.005 mg/L 1 05/07/20 TH < 0.010 0.010 1 05/07/20 SW6010D Selenium mg/L TH < 0.001 0.001 mg/L 1 05/07/20 RS SW7010 Thallium 0.012 0.002 05/07/20 TH SW6010D Vanadium mg/L 1 Zinc 0.032 0.004 mg/L 05/07/20 TH SW6010D Completed /BF 0.45um Filter Filtration Completed 05/07/20 RA/RA SW7470A Mercury Digestion Completed 05/06/20 AG/BF **Total Metals Digestion**

Client ID: MW-2 25UM

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

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

Phyllis Shiller, Laboratory Director

May 14, 2020

Reviewed and Released by: Phyllis Shiller, Laboratory Director



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

Analysis Report

May 14, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Sample Information Custody Information Date <u>Time</u>

GROUND WATER Collected by: 05/05/20 Matrix:

Received by: **BLCOMPANIES** В 05/06/20 **Location Code:** 15:46

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: **Laboratory Data** SDG ID: GCF90108

Phoenix ID: CF90198

Project ID: 16D3253 Client ID: MW-1 1.2UM

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	05/07/20	TH	SW6010D
Barium	0.054	0.002	mg/L	1	05/07/20	TH	SW6010D
Beryllium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Chromium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Mercury	< 0.0002	0.0002	mg/L	1	05/08/20	RS	SW7470A
Nickel	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Lead	< 0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Antimony	< 0.005	0.005	mg/L	1	05/07/20	TH	SW6010D
Selenium	< 0.010	0.010	mg/L	1	05/07/20	TH	SW6010D
Thallium	< 0.001	0.001	mg/L	1	05/07/20	RS	SW7010
Vanadium	0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Zinc	< 0.004	0.004	mg/L	1	05/07/20	TH	SW6010D
Filtration	Completed					/BF	0.45um Filter
Mercury Digestion	Completed				05/07/20	RA/RA	SW7470A
Total Metals Digestion	Completed				05/06/20	AG/BF	:

Client ID: MW-1 1.2UM

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

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

Phyllis Shiller, Laboratory Director

May 14, 2020

Reviewed and Released by: Phyllis Shiller, Laboratory Director



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

Analysis Report

May 14, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: GROUND WATER Collected by: 05/05/20

Location Code: BLCOMPANIES Received by: B 05/06/20 15:46

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF90108

Phoenix ID: CF90199

Project ID: 16D3253 Client ID: MW-2 1.2UM

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	05/07/20	TH	SW6010D
Barium	0.059	0.002	mg/L	1	05/07/20	TH	SW6010D
Beryllium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Chromium	< 0.001	0.001	mg/L	1	05/07/20	TH	SW6010D
Mercury	< 0.0002	0.0002	mg/L	1	05/08/20	RS	SW7470A
Nickel	0.005	0.001	mg/L	1	05/07/20	TH	SW6010D
Lead	< 0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Antimony	< 0.005	0.005	mg/L	1	05/07/20	TH	SW6010D
Selenium	< 0.010	0.010	mg/L	1	05/07/20	TH	SW6010D
Thallium	< 0.001	0.001	mg/L	1	05/07/20	RS	SW7010
Vanadium	< 0.002	0.002	mg/L	1	05/07/20	TH	SW6010D
Zinc	0.010	0.004	mg/L	1	05/07/20	TH	SW6010D
Filtration	Completed					/BF	0.45um Filter
Mercury Digestion	Completed				05/07/20	RA/RA	SW7470A
Total Metals Digestion	Completed				05/06/20	AG/BF	•

Client ID: MW-2 1.2UM

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

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

Phyllis Shiller, Laboratory Director

May 14, 2020

Reviewed and Released by: Phyllis Shiller, Laboratory Director



Tuesday, March 24, 2020

Attn: Matthew Gonsalves BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Project ID: 16D3253 SDG ID: GCF51973

Sample ID#s: CF51973 - CF51979

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

Sincerely yours,

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



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



SDG Comments

March 24, 2020

SDG I.D.: GCF51973

8260 Analysis:

1,2-Dibromoethane doesn't meet GW-1 criteria, this compound is analyzed by GC/FID to achieve this criteria.

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

Phoenix reporting levels may exceed those referenced in the CAM protocol. Please refer to criteria sheet for comparisons to requested MCP standards.



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

Sample Id Cross Reference

March 24, 2020

SDG I.D.: GCF51973

Project ID: 16D3253

Client Id	Lab Id	Matrix
MW-1	CF51973	WATER
MW-2	CF51974	WATER
MW-101	CF51975	WATER
MW-102	CF51976	WATER
MW-103	CF51977	WATER
MW-104	CF51978	WATER
TRIP BLANK	CF51979	WATER



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51973

Project ID: 16D3253 Client ID: MW-1

RL/

Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Volatiles							_
	ND	1.0	//	4	03/17/20	МН	SW8260C
1,1,1,2-Tetrachloroethane	ND ND	1.0	ug/L	1	03/17/20	МН	SW8260C SW8260C
1,1,1-Trichloroethane			ug/L	1			
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2-Hexanone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
2-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
4-Methyl-2-pentanone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
wellyi-2-pentanone	ND	5.0	ug/ L	1	00/11/20	IVII I	01102000

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
	ND	25	ug/L	1	03/17/20	MH	SW8260C
Acetone Acrylonitrile	ND	1.0	ug/∟ ug/L	1	03/17/20	MH	SW8260C SW8260C
Benzene	ND	0.70	ug/L	1	03/17/20	MH	SW8260C
Bromobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,2-Dichloroethene	1.7	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Dibromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Dibromomethane Dishlaradifluoremethane					03/17/20		
Dichlorodifluoromethane	ND ND	1.0 1.0	ug/L	1 1	03/17/20	MH MH	SW8260C SW8260C
Ethylbenzene	ND	0.40	ug/L	1	03/17/20	МН	SW8260C SW8260C
Hexachlorobutadiene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Isopropylbenzene			ug/L				
m&p-Xylene	ND	1.0 5.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C SW8260C
Methyl ethyl ketone	ND		ug/L	1	03/17/20	MH MH	
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1			SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Naphthalene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Propylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
o-Xylene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
p-Isopropyltoluene	ND ND	1.0 1.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C SW8260C
sec-Butylbenzene			ug/L	1		MH MH	
Styrene	ND	1.0	ug/L	•	03/17/20		SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrachloroethene	ND	1.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1		MH	SW8260C
Toluene	ND	1.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C
Total Xylenes	ND	1.0	ug/L	1		MH	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Trichloroethene	1.7	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorotrifluoroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Vinyl chloride	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
QA/QC Surrogates	05		07	4	00/47/00	N 41 1	70 420 0/
% 1,2-dichlorobenzene-d4	95		%	1	03/17/20	MH	70 - 130 %
% Bromofluorobenzene	93		%	1	03/17/20	MH	70 - 130 %
% Dibromofluoromethane	84		%	1	03/17/20	МН	70 - 130 %

Client ID: MW-1

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
% Toluene-d8	95		%	1	03/17/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	03/17/20	МН	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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:

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: SDG ID: GCF51973 Phoenix ID: CF51974

Project ID: 16D3253 Client ID: MW-2

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Arsenic	< 0.004	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Barium	0.106	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Beryllium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Cadmium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Chromium	0.002	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury	< 0.0002	0.0002	mg/L	1	03/18/20	RS	SW7470/E245.1
Nickel	0.004	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Lead	< 0.002	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Antimony	< 0.005	0.005	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Selenium	< 0.010	0.010	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Thallium	< 0.0005	0.0005	mg/L	5	03/19/20	CPP	SW6020B/E200.8
Vanadium	0.003	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Zinc	0.017	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury Digestion	Completed				03/18/20	3/RA/LS/	RSW7470/245.1
EPH Extraction	Completed				03/18/20	SB/SB	SW3510C
MA Petroleum Hydrocarbon (EPH)	Completed				03/17/20		MADEP EPH-04
Semi-Volatile Extraction	Completed				03/17/20	P/AK	SW3520C
Total Metals Digestion	Completed				03/17/20	AG	
Total Metals Digestion MS	Completed				03/18/20	AG	
MA Petroleum Hydrocarbon (VPH)	Completed				03/18/20	RM	MADEP VPH04
<u>Volatiles</u>							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1-Dichloroethene 1,1-Dichloropropene	ND	1.0	ug/∟ ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2,2-Dichloropropane2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
2-Hexanone					03/17/20		
2-Isopropyltoluene	ND ND	1.0 1.0	ug/L ug/L	1 1	03/17/20	MH MH	SW8260C SW8260C
4-Chlorotoluene	ND	5.0		1	03/17/20	МН	SW8260C SW8260C
4-Methyl-2-pentanone	ND	25	ug/L ug/L	1	03/17/20	MH	SW8260C
Acetone							
Acrylonitrile	ND ND	1.0 0.70	ug/L	1	03/17/20 03/17/20	MH	SW8260C SW8260C
Benzene			ug/L	1	03/17/20	MH MH	
Bromobenzene	ND	1.0	ug/L	1			SW8260C
Bromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND ND	1.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C SW8260C
Chlorobenzene	ND	1.0	ug/L	1		MH MH	SW8260C SW8260C
Chloroethane		1.0	ug/L	•	03/17/20		
Chloroform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C SW8260C
cis-1,2-Dichloroethene	3.7 ND	1.0	ug/L	1	03/17/20	MH	
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Dibromochloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Dibromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Dichlorodifluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Ethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Isopropylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
m&p-Xylene	ND	1.0	ug/L	1 4	03/17/20	MH	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Naphthalene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
n-Propylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
o-Xylene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
p-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
sec-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Styrene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrachloroethene	4.5	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	03/17/20	MH	SW8260C
Toluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Total Xylenes	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Trichloroethene	20	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorotrifluoroethane	79	5.0	ug/L	5	03/18/20	MH	SW8260C
Vinyl chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
QA/QC Surrogates							
% 1,2-dichlorobenzene-d4	97		%	1	03/17/20	MH	70 - 130 %
% Bromofluorobenzene	93		%	1	03/17/20	MH	70 - 130 %
% Dibromofluoromethane	88		%	1	03/17/20	MH	70 - 130 %
% Toluene-d8	103		%	1	03/17/20	MH	70 - 130 %
% 1,2-dichlorobenzene-d4 (5x)	95		%	5	03/18/20	MH	70 - 130 %
% Bromofluorobenzene (5x)	91		%	5	03/18/20	MH	70 - 130 %
% Dibromofluoromethane (5x)	88		%	5	03/18/20	MH	70 - 130 %
% Toluene-d8 (5x)	96		%	5	03/18/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	03/17/20	MH	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)
<u>Semivolatiles</u>							
1,2,4,5-Tetrachlorobenzene	ND	3.6	ug/L	1	03/20/20	WB	SW8270D
1,2,4-Trichlorobenzene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
1,2-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
1,2-Diphenylhydrazine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
1,3-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
1,4-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
2,4,5-Trichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4,6-Trichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dimethylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dinitrophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dinitrotoluene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2,6-Dinitrotoluene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Chloronaphthalene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Chlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
2-Methylphenol (o-cresol)	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Nitrophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	10	ug/L	1	03/20/20	WB	SW8270D
3,3'-Dichlorobenzidine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
3-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Bromophenyl phenyl ether	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Chloro-3-methylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Chloroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Nitrophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Acetophenone	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Aniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Benzidine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Benzoic acid	ND	51	ug/L	1	03/20/20	WB	SW8270D
Benzyl butyl phthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethyl)ether	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Carbazole	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Dibenzofuran	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Diethyl phthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Dimethylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Di-n-butylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Di-n-octylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Hexachloroethane	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Isophorone	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodiphenylamine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Pentachloronitrobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
Phenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
QA/QC Surrogates	110	1.0	ug, L	•	00/20/20	****	01102103
% 2,4,6-Tribromophenol	71		%	1	03/20/20	WB	15 - 110 %
% 2-Fluorobiphenyl	71		%	1	03/20/20	WB	30 - 130 %
% 2-Fluorophenol	54		%	1	03/20/20	WB	15 - 110 %
% Nitrobenzene-d5	68		%	1	03/20/20	WB	30 - 130 %
% Phenol-d5	11		%	1	03/20/20	WB	15 - 110 % ³
% Terphenyl-d14	47		%	1	03/20/20	WB	30 - 130 %
	71		70	•	00/20/20	***	00 100 /0
Semivolatiles (SIM)		0.15	_		00/:0/0-		014/00752 (01/5)
2-Methylnaphthalene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthylene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Anthracene	ND	0.09	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benz(a)anthracene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(a)pyrene	ND	0.20	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(b)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)

Client ID. WW-2		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Benzo(ghi)perylene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(k)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Chrysene	ND	0.05	ug/L	1	03/18/20	WB	SW8270D (SIM)
Dibenz(a,h)anthracene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)
Fluoranthene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Fluorene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorobenzene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorobutadiene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorocyclopentadiene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Indeno(1,2,3-cd)pyrene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Naphthalene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Nitrobenzene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
N-Nitrosodimethylamine	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pentachlorophenol	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Phenanthrene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pyrene	ND	0.07	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pyridine	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
QA/QC Surrogates							
% 2,4,6-Tribromophenol	84		%	1	03/18/20	WB	15 - 110 %
% 2-Fluorobiphenyl	76		%	1	03/18/20	WB	40 - 140 %
% 2-Fluorophenol	69		%	1	03/18/20	WB	15 - 110 %
% Nitrobenzene-d5	72		%	1	03/18/20	WB	40 - 140 %
% Phenol-d5	77		%	1	03/18/20	WB	15 - 110 %
% Terphenyl-d14	62		%	1	03/18/20	WB	40 - 140 %
MA EPH Aliphatic/Aroma	atic Rang	<u>jes</u>					
C11-C22 Aromatic Hydrocarbons 1,2*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C11-C22 Aromatic Hydrocarbons Unadj	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C19-C36 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C9-C18 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
Total TPH 1,2*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
QA/QC Surrogates							
% 1-chlorooctadecane (aliphatic)	58		%	1	03/18/20	AW	40 - 140 %
% 2-Bromonaphthalene (Fractionation)	112		%	1	03/18/20	AW	40 - 140 %
% 2-Fluorobiphenyl (Fractionation)	109		%	1	03/18/20	AW	40 - 140 %
% o-terphenyl (aromatic)	88		%	1	03/18/20	AW	40 - 140 %
MA Volatile Petroleum H	ydrocark	ons (VP	<u>H)</u>				
Unadjusted C5-C8 Aliphatics (*1)	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
Unadjusted C9-C12 Aliphatics (*1)	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C5-C8 Aliphatic Hydrocarbons *1,2	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C9-C12 Aliphatic Hydrocarbons *1,3	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C9-C10 Aromatic Hydrocarbons *1	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Ethyl Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
MTBE	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Naphthalene	ND	5.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Toluene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
m,p-Xylenes	ND	2.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
o-Xylene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004

Client ID: MW-2

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
QA/QC Surrogates							
% 2,5-Dibromotoluene (FID)	95		%	1	03/18/20	RM	70 - 130 %
% 2,5-Dibromotoluene (PID)	92		%	1	03/18/20	RM	70 - 130 %

^{3 =} This parameter exceeds laboratory specified limits.

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level 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.

MAEPH:

- 1* Hydrocarbon range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
- 2* C11-C12 Aromatic Hydrocarbons exclude the concentration of Target PAH analytes eluting in that range.

VPH:

- *1 Range data exclude conc.s of any surrogate(s) and/or Int. std.s eluting in that range.
- *2 C5-C8 and C9-C12 Aliphatic exclude the conc. of Target Analytes in that range.
- *3 C9-C12 Aliphatic also exclude C9-C10 Aromatic Hydrocarbon

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

Semi-Volatile Comment:

Poor surrogate recovery was observed for one acid and/or one base surrogate. The other surrogates associated with this sample were within QA/QC criteria. No significant bias suspected.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51975

Project ID: 16D3253 Client ID: MW-101

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Arsenic	0.007	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Barium	0.099	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Beryllium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Cadmium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Chromium	0.012	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury	< 0.0002	0.0002	mg/L	1	03/18/20	RS	SW7470/E245.1
Nickel	0.008	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Lead	0.009	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Antimony	< 0.005	0.005	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Selenium	< 0.010	0.010	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Thallium	< 0.0005	0.0005	mg/L	5	03/19/20	CPP	SW6020B/E200.8
Vanadium	0.009	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Zinc	0.017	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury Digestion	Completed				03/18/20	3/RA/LS/	RSW7470/245.1
EPH Extraction	Completed				03/18/20	SB/SB	SW3510C
MA Petroleum Hydrocarbon (EPH)	Completed				03/17/20		MADEP EPH-04
Semi-Volatile Extraction	Completed				03/17/20	P/AK	SW3520C
Total Metals Digestion	Completed				03/17/20	AG	
Total Metals Digestion MS	Completed				03/18/20	AG	
MA Petroleum Hydrocarbon (VPH)	Completed				03/18/20	RM	MADEP VPH04
<u>Volatiles</u>							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1-Dichloroethene 1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C SW8260C
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichloropropane 1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	8.1	1.0	ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2,4-Trimethylbenzene	ND	1.0		1	03/17/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2-Dibromoethane	ND	1.0	ug/L ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2-Dichlorobenzene	ND	0.60	ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2-Dichloroethane							
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3,5-Trimethylbenzene	3.3 ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2-Hexanone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
2-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
4-Methyl-2-pentanone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
Acetone	ND	25	ug/L	1	03/17/20	МН	SW8260C
Acrylonitrile	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Benzene	ND	0.70	ug/L	1	03/17/20	МН	SW8260C
Bromobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,2-Dichloroethene	6.4	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Dibromochloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Dibromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Dichlorodifluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Ethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Isopropylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
m&p-Xylene	4.3	1.0	ug/L	1	03/17/20	МН	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Naphthalene	2.7	1.0	ug/L	1	03/17/20	МН	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
•			-				

December 1	D II	RL/	11.5.	Dil die	Data/Time	_	Defense
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
n-Propylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
o-Xylene	2.9	1.0	ug/L	1	03/17/20	МН	SW8260C
p-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
sec-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Styrene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrachloroethene	7.0	1.0	ug/L	1	03/17/20	МН	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	03/17/20	МН	SW8260C
Toluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Total Xylenes	7.2	1.0	ug/L	1	03/17/20	МН	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	МН	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Trichloroethene	44	5.0	ug/L	5	03/18/20	MH	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorotrifluoroethane	170	20	ug/L	20	03/19/20	MH	SW8260C
Vinyl chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
QA/QC Surrogates							
% 1,2-dichlorobenzene-d4	96		%	1	03/17/20	MH	70 - 130 %
% Bromofluorobenzene	94		%	1	03/17/20	MH	70 - 130 %
% Dibromofluoromethane	85		%	1	03/17/20	MH	70 - 130 %
% Toluene-d8	107		%	1	03/17/20	MH	70 - 130 %
% 1,2-dichlorobenzene-d4 (5x)	96		%	5	03/18/20	MH	70 - 130 %
% Bromofluorobenzene (5x)	93		%	5	03/18/20	MH	70 - 130 %
% Dibromofluoromethane (5x)	89		%	5	03/18/20	МН	70 - 130 %
% Toluene-d8 (5x)	98		%	5	03/18/20	МН	70 - 130 %
% 1,2-dichlorobenzene-d4 (20x)	101		%	20	03/19/20	МН	70 - 130 %
% Bromofluorobenzene (20x)	93		%	20	03/19/20	МН	70 - 130 %
% Dibromofluoromethane (20x)	101		%	20	03/19/20	МН	70 - 130 %
% Toluene-d8 (20x)	99		%	20	03/19/20	МН	70 - 130 %
O							
Oxygenates & Dioxane	ND	50	//	4	02/47/00	N 41 1	CM00000 (OVV)
1,4-Dioxane	ND	50	ug/L	1	03/17/20	MH	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)
<u>Semivolatiles</u>							
1,2,4,5-Tetrachlorobenzene	ND	3.6	ug/L	1	03/20/20	WB	SW8270D
1,2,4-Trichlorobenzene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
1,2-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
1,2-Diphenylhydrazine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
1,3-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
1,4-Dichlorobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
2,4,5-Trichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4,6-Trichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dichlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2,4-Dimethylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
					03/20/20		SW8270D SW8270D
2,4-Dinitrophenol	ND	1.0	ug/L	1	03/20/20	WB	34405100

Chefft ID. WWV-101		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
2,4-Dinitrotoluene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2,6-Dinitrotoluene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Chloronaphthalene	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Chlorophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2-Methylphenol (o-cresol)	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
2-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
2-Nitrophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	10	ug/L	1	03/20/20	WB	SW8270D
3,3'-Dichlorobenzidine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
3-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Bromophenyl phenyl ether	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Chloro-3-methylphenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Chloroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
4-Nitroaniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
4-Nitrophenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Acetophenone	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Aniline	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Benzidine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Benzoic acid	ND	51	ug/L	1	03/20/20	WB	SW8270D
Benzyl butyl phthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethyl)ether	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Carbazole	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Dibenzofuran	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Diethyl phthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Dimethylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Di-n-butylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Di-n-octylphthalate	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Hexachloroethane	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
Isophorone	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodiphenylamine	ND	5.1	ug/L	1	03/20/20	WB	SW8270D
Pentachloronitrobenzene	ND	2.6	ug/L	1	03/20/20	WB	SW8270D
Phenol	ND	1.0	ug/L	1	03/20/20	WB	SW8270D
QA/QC Surrogates			-9-		55,25,25		
% 2,4,6-Tribromophenol	68		%	1	03/20/20	WB	15 - 110 %
% 2-Fluorobiphenyl	69		%	1	03/20/20	WB	30 - 130 %
% 2-Fluorophenol	49		%	1	03/20/20	WB	15 - 110 %
% Nitrobenzene-d5	73		%	1	03/20/20	WB	30 - 130 %
% Phenol-d5	40		%	1	03/20/20	WB	15 - 110 %
% Terphenyl-d14	42		%	1	03/20/20	WB	30 - 130 %
Semivolatiles (SIM)							
2-Methylnaphthalene	1.2	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthylene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
. Toonapharylene	110	5.10	49/ ∟	'	33, 10,20	***	211321 0D (OIIVI)

Davamatar	Decult	RL/	l loite	Dilution	Data/Times	D	Deference	
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference	
Anthracene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Benz(a)anthracene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Benzo(a)pyrene	ND	0.21	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Benzo(b)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Benzo(ghi)perylene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Benzo(k)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Chrysene	ND	0.05	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Dibenz(a,h)anthracene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Fluoranthene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Fluorene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Hexachlorobenzene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Hexachlorobutadiene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Hexachlorocyclopentadiene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Indeno(1,2,3-cd)pyrene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Naphthalene	1.6	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Nitrobenzene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
N-Nitrosodimethylamine	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Pentachlorophenol	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Phenanthrene	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Pyrene	0.12	0.07	ug/L	1	03/18/20	WB	SW8270D (SIM)	
Pyridine	ND	0.52	ug/L	1	03/18/20	WB	SW8270D (SIM)	
QA/QC Surrogates								
% 2,4,6-Tribromophenol	85		%	1	03/18/20	WB	15 - 110 %	
% 2-Fluorobiphenyl	64		%	1	03/18/20	WB	40 - 140 %	
% 2-Fluorophenol	45		%	1	03/18/20	WB	15 - 110 %	
% Nitrobenzene-d5	67		%	1	03/18/20	WB	40 - 140 %	
% Phenol-d5	57		%	1	03/18/20	WB	15 - 110 %	
% Terphenyl-d14	20		%	1	03/18/20	WB	40 - 140 %	3
	d'a Bassa							
MA EPH Aliphatic/Aroma			,		00/00/00	****		
C11-C22 Aromatic Hydrocarbons 1,2*	ND	200	ug/L	1	03/20/20	AW	MAEPH 5/2004	
C11-C22 Aromatic Hydrocarbons Unadj	ND	200	ug/L	1	03/20/20	AW	MAEPH 5/2004	
C19-C36 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/20/20	AW	MAEPH 5/2004	
C9-C18 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/20/20	AW	MAEPH 5/2004	
Total TPH 1,2*	ND	200	ug/L	1	03/20/20	AW	MAEPH 5/2004	
QA/QC Surrogates								
% 1-chlorooctadecane (aliphatic)	62		%	1	03/20/20	AW	40 - 140 %	
% 2-Bromonaphthalene (Fractionation)	106		%	1	03/20/20	AW	40 - 140 %	
% 2-Fluorobiphenyl (Fractionation)	87		%	1	03/20/20	AW	40 - 140 %	
% o-terphenyl (aromatic)	66		%	1	03/20/20	AW	40 - 140 %	
MA Volatile Petroleum H	ydrocark	ons (VP	<u>PH)</u>					
Unadjusted C5-C8 Aliphatics (*1)	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004	
Unadjusted C9-C12 Aliphatics (*1)	150	100	ug/L	1	03/18/20	RM	MA VPH 5/2004	
C5-C8 Aliphatic Hydrocarbons *1,2	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004	
C9-C12 Aliphatic Hydrocarbons *1,3	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004	
C9-C10 Aromatic Hydrocarbons *1	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004	
Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004	
Ethyl Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004	
MTBE	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004	
-		*	- 3					

Client ID: MW-101

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Naphthalene	ND	5.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Toluene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
m,p-Xylenes	4.1	2.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
o-Xylene	3.1	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
QA/QC Surrogates							
% 2,5-Dibromotoluene (FID)	99		%	1	03/18/20	RM	70 - 130 %
% 2,5-Dibromotoluene (PID)	93		%	1	03/18/20	RM	70 - 130 %

^{3 =} This parameter exceeds laboratory specified limits.

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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.

MAEPH:

- 1* Hydrocarbon range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
- 2* C11-C12 Aromatic Hydrocarbons exclude the concentration of Target PAH analytes eluting in that range.

VPH:

- *1 Range data exclude conc.s of any surrogate(s) and/or Int. std.s eluting in that range.
- *2 C5-C8 and C9-C12 Aliphatic exclude the conc. of Target Analytes in that range.
- *3 C9-C12 Aliphatic also exclude C9-C10 Aromatic Hydrocarbon

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

Semi-Volatile Comment:

Poor surrogate recovery was observed for one acid and/or one base surrogate. The other surrogates associated with this sample were within QA/QC criteria. No significant bias suspected.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51976

Project ID: 16D3253 Client ID: MW-102

RL/

Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Volatiles							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2-Hexanone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
2-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
4-Methyl-2-pentanone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
	ND						SW8260C
Acetone	ND	25 1.0	ug/L	1 1	03/17/20 03/17/20	MH MH	SW8260C SW8260C
Acrylonitrile	ND	0.70	ug/L	1	03/17/20	МН	SW8260C SW8260C
Benzene Bromobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C SW8260C
Bromochloromethane	ND	1.0	ug/L ug/L	1	03/17/20	МН	SW8260C SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,2-Dichloroethene	5.4	1.0	ug/L	1	03/17/20	MH	SW8260C
,	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
cis-1,3-Dichloropropene Dibromochloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Dibromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Dichlorodifluoromethane	2.6	1.0	ug/L	1	03/17/20	MH	SW8260C
Ethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Isopropylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
m&p-Xylene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Naphthalene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Propylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
o-Xylene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
p-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
sec-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Styrene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrachloroethene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	03/17/20	MH	SW8260C
Toluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Total Xylenes	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	МН	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
Trichloroethene	5.8	1.0	ug/L	1	03/17/20	МН	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Trichlorotrifluoroethane	2.4	1.0	ug/L	1	03/17/20	MH	SW8260C
Vinyl chloride	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
QA/QC Surrogates		-	. –				
% 1,2-dichlorobenzene-d4	96		%	1	03/17/20	МН	70 - 130 %
% Bromofluorobenzene	93		%	1	03/17/20	МН	70 - 130 %
% Dibromofluoromethane	86		%	1	03/17/20	МН	70 - 130 %
					-		

Client ID: MW-102

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
% Toluene-d8	98		%	1	03/17/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	03/17/20	MH	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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:

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51977

Project ID: 16D3253 Client ID: MW-103

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Arsenic	< 0.004	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Barium	0.130	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Beryllium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Cadmium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Chromium	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury	< 0.0002	0.0002	mg/L	1	03/18/20	RS	SW7470/E245.1
Nickel	< 0.001	0.001	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Lead	< 0.002	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Antimony	< 0.005	0.005	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Selenium	< 0.010	0.010	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Thallium	< 0.0005	0.0005	mg/L	5	03/19/20	CPP	SW6020B/E200.8
Vanadium	< 0.002	0.002	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Zinc	< 0.004	0.004	mg/L	1	03/18/20	CPP	SW6010D/E200.7
Mercury Digestion	Completed				03/18/20	3/RA/LS/	RSW7470/245.1
EPH Extraction	Completed				03/18/20	SB/SB	SW3510C
MA Petroleum Hydrocarbon (EPH)	Completed				03/17/20		MADEP EPH-04
Semi-Volatile Extraction	Completed				03/17/20	P/AK	SW3520C
Total Metals Digestion	Completed				03/17/20	AG	
Total Metals Digestion MS	Completed				03/18/20	AG	
MA Petroleum Hydrocarbon (VPH)	Completed				03/18/20	RM	MADEP VPH04
<u>Volatiles</u>							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C

Project ID: 16D3253 Phoenix I.D.: CF51977 Client ID: MW-103

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1-Dichloroethene 1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	MH	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2,2-Dichloropropane2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
2-Hexanone					03/17/20		
2-Isopropyltoluene	ND ND	1.0 1.0	ug/L	1	03/17/20	MH MH	SW8260C SW8260C
4-Chlorotoluene			ug/L	1			
4-Methyl-2-pentanone	ND	5.0	ug/L	1	03/17/20 03/17/20	MH	SW8260C
Acetone	ND	25	ug/L	1		MH	SW8260C
Acrylonitrile	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Benzene	ND	0.70	ug/L	1	03/17/20	MH	SW8260C
Bromobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,2-Dichloroethene	30 ND	1.0	ug/L	1	03/17/20	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Dibromochloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Dibromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Dichlorodifluoromethane	7.1	1.0	ug/L	1	03/17/20	MH	SW8260C
Ethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
Isopropylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
m&p-Xylene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Naphthalene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C

Client ID: MW-103 RL/ Parameter Result **PQL** Units Dilution Date/Time Βv Reference ND n-Propylbenzene 1.0 ug/L 1 03/17/20 MH SW8260C 1 o-Xylene ND ug/L 03/17/20 MH SW8260C 1.0 p-Isopropyltoluene ND 1.0 ug/L 1 03/17/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 03/17/20 MH SW8260C ND 1.0 ug/L 1 MH SW8260C Styrene 03/17/20 1 tert-Butylbenzene ND 1.0 ug/L 03/17/20 MH SW8260C 1 5.4 1.0 ug/L SW8260C Tetrachloroethene 03/17/20 MH ND 1 Tetrahydrofuran (THF) 2.5 ug/L 03/17/20 MH SW8260C ug/L ND 1.0 1 MH SW8260C Toluene 03/17/20 1 SW8260C Total Xylenes ND 1.0 ug/L 03/17/20 MH ND 1.0 ug/L 1 03/17/20 MH SW8260C trans-1,2-Dichloroethene trans-1,3-Dichloropropene ND 0.40 ug/L 1 03/17/20 MH SW8260C 5.0 1 trans-1,4-dichloro-2-butene ND ug/L 03/17/20 MH SW8260C 5 34 5.0 ug/L 03/18/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 03/17/20 MH SW8260C Trichlorofluoromethane Trichlorotrifluoroethane 84 5.0 ug/L 5 03/18/20 MH SW8260C ND 1.0 ug/L 1 03/17/20 MH SW8260C Vinyl chloride **QA/QC Surrogates** 97 % 1 03/17/20 70 - 130 % % 1.2-dichlorobenzene-d4 MH 1 94 % 03/17/20 MH 70 - 130 % % Bromofluorobenzene 85 % 1 03/17/20 MH 70 - 130 % % Dibromofluoromethane % Toluene-d8 107 % 1 03/17/20 MH 70 - 130 % % 1,2-dichlorobenzene-d4 (5x) 99 % 5 03/18/20 MH 70 - 130 % 5 % Bromofluorobenzene (5x) 93 % 03/18/20 MH 70 - 130 % 100 % 5 03/18/20 МН 70 - 130 % % Dibromofluoromethane (5x) 97 % 5 03/18/20 70 - 130 % % Toluene-d8 (5x) MH Oxygenates & Dioxane 1,4-Dioxane ND 50 ug/L 1 03/17/20 МН SW8260C (OXY) ND 1.0 ug/L 1 03/17/20 МН SW8260C (OXY) Diethyl ether ND ug/L 1 03/17/20 SW8260C (OXY) Di-isopropyl ether 1.0 MH Ethyl tert-butyl ether 03/17/20 ND 1.0 ug/L 1 MH SW8260C (OXY) ND 1 03/17/20 MH SW8260C (OXY) tert-amyl methyl ether 1.0 ug/L **Semivolatiles** 1 03/20/20 1,2,4,5-Tetrachlorobenzene ND 3.9 ug/L WB SW8270D 03/20/20 SW8270D ND 5.5 ug/L 1 WB 1,2,4-Trichlorobenzene ND 2.8 ug/L 1 03/20/20 WB SW8270D 1,2-Dichlorobenzene SW8270D ND 5.5 ug/L 1 03/20/20 WB 1,2-Diphenylhydrazine 1,3-Dichlorobenzene ND 2.8 ug/L 1 03/20/20 WB SW8270D 2.8 1 SW8270D ND ug/L 03/20/20 WB 1,4-Dichlorobenzene 03/20/20 SW8270D 2,4,5-Trichlorophenol ND 1.1 ug/L 1 WB

ug/L

ug/L

ug/L

ug/L

ug/L

ug/L

ug/L

ug/L

1

1

1

1

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SW8270D

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SW8270D

SW8270D

SW8270D

SW8270D

SW8270D

SW8270D

ND

ND

ND

ND

ND

ND

ND

ND

1.1

1.1

1.1

1.1

5.5

5.5

5.5

1.1

2,4,6-Trichlorophenol

2,4-Dichlorophenol

2,4-Dimethylphenol

2,4-Dinitrophenol

2,4-Dinitrotoluene

2,6-Dinitrotoluene

2-Chlorophenol

2-Chloronaphthalene

Project ID: 16D3253 Phoenix I.D.: CF51977 Client ID: MW-103

Client ID. MW-103		DL /					
Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
2-Methylphenol (o-cresol)	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
2-Nitroaniline	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
2-Nitrophenol	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	11	ug/L	1	03/20/20	WB	SW8270D
3,3'-Dichlorobenzidine	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
3-Nitroaniline	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
4-Bromophenyl phenyl ether	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
4-Chloro-3-methylphenol	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
4-Chloroaniline	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
4-Nitroaniline	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
4-Nitrophenol	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
Acetophenone	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Aniline	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Benzidine	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Benzoic acid	ND	55	ug/L	1	03/20/20	WB	SW8270D
Benzyl butyl phthalate	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroethyl)ether	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
Carbazole	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Dibenzofuran	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Diethyl phthalate	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Dimethylphthalate	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Di-n-butylphthalate	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Di-n-octylphthalate	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Hexachloroethane	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
Isophorone	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
N-Nitrosodiphenylamine	ND	5.5	ug/L	1	03/20/20	WB	SW8270D
Pentachloronitrobenzene	ND	2.8	ug/L	1	03/20/20	WB	SW8270D
Phenol	ND	1.1	ug/L	1	03/20/20	WB	SW8270D
QA/QC Surrogates							
% 2,4,6-Tribromophenol	85		%	1	03/20/20	WB	15 - 110 %
% 2-Fluorobiphenyl	76		%	1	03/20/20	WB	30 - 130 %
% 2-Fluorophenol	51		%	1	03/20/20	WB	15 - 110 %
% Nitrobenzene-d5	72		%	1	03/20/20	WB	30 - 130 %
% Phenol-d5	33		%	1	03/20/20	WB	15 - 110 %
% Terphenyl-d14	81		%	1	03/20/20	WB	30 - 130 %
Semivolatiles (SIM)							
2-Methylnaphthalene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Acenaphthylene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Anthracene	ND	0.09	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benz(a)anthracene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(a)pyrene	ND	0.19	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(b)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)

Parameter Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Benzo(ghi)perylene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)
Benzo(k)fluoranthene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Chrysene	ND	0.05	ug/L	1	03/18/20	WB	SW8270D (SIM)
Dibenz(a,h)anthracene	ND	0.02	ug/L	1	03/18/20	WB	SW8270D (SIM)
Fluoranthene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Fluorene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorobenzene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorobutadiene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Hexachlorocyclopentadiene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Indeno(1,2,3-cd)pyrene	ND	0.10	ug/L	1	03/18/20	WB	SW8270D (SIM)
Naphthalene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Nitrobenzene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
N-Nitrosodimethylamine	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pentachlorophenol	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Phenanthrene	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pyrene	ND	0.07	ug/L	1	03/18/20	WB	SW8270D (SIM)
Pyridine	ND	0.49	ug/L	1	03/18/20	WB	SW8270D (SIM)
QA/QC Surrogates							
% 2,4,6-Tribromophenol	85		%	1	03/18/20	WB	15 - 110 %
% 2-Fluorobiphenyl	73		%	1	03/18/20	WB	40 - 140 %
% 2-Fluorophenol	59		%	1	03/18/20	WB	15 - 110 %
% Nitrobenzene-d5	66		%	1	03/18/20	WB	40 - 140 %
% Phenol-d5	70		%	1	03/18/20	WB	15 - 110 %
% Terphenyl-d14	67		%	1	03/18/20	WB	40 - 140 %
MA EPH Aliphatic/Aroma	atic Rang	<u>jes</u>					
C11-C22 Aromatic Hydrocarbons 1,2*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C11-C22 Aromatic Hydrocarbons Unadj	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C19-C36 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
C9-C18 Aliphatic Hydrocarbons 1*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
Total TPH 1,2*	ND	200	ug/L	1	03/18/20	AW	MAEPH 5/2004
QA/QC Surrogates							
% 1-chlorooctadecane (aliphatic)	64		%	1	03/18/20	AW	40 - 140 %
% 2-Bromonaphthalene (Fractionation)	111		%	1	03/18/20	AW	40 - 140 %
% 2-Fluorobiphenyl (Fractionation)	108		%	1	03/18/20	AW	40 - 140 %
% o-terphenyl (aromatic)	86		%	1	03/18/20	AW	40 - 140 %
MA Volatile Petroleum H	ydrocarl	ons (VPI	<u>1)</u>				
Unadjusted C5-C8 Aliphatics (*1)	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
Unadjusted C9-C12 Aliphatics (*1)	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C5-C8 Aliphatic Hydrocarbons *1,2	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C9-C12 Aliphatic Hydrocarbons *1,3	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
C9-C10 Aromatic Hydrocarbons *1	ND	100	ug/L	1	03/18/20	RM	MA VPH 5/2004
Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Ethyl Benzene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
MTBE	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Naphthalene	ND	5.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
Toluene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
m,p-Xylenes	ND	2.0	ug/L	1	03/18/20	RM	MA VPH 5/2004
o-Xylene	ND	1.0	ug/L	1	03/18/20	RM	MA VPH 5/2004

Client ID: MW-103

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
QA/QC Surrogates % 2,5-Dibromotoluene (FID) % 2,5-Dibromotoluene (PID)	90 88		% %	1 1	03/18/20 03/18/20	RM RM	70 - 130 % 70 - 130 %

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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.

MAEPH:

- 1* Hydrocarbon range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
- 2* C11-C12 Aromatic Hydrocarbons exclude the concentration of Target PAH analytes eluting in that range.

VPH:

- *1 Range data exclude conc.s of any surrogate(s) and/or Int. std.s eluting in that range.
- *2 C5-C8 and C9-C12 Aliphatic exclude the conc. of Target Analytes in that range.
- *3 C9-C12 Aliphatic also exclude C9-C10 Aromatic Hydrocarbon

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51978

Project ID: 16D3253 Client ID: MW-104

RL/

Parameter	meter Result PQL Units	Dilution	Date/Time	Ву	Reference		
Volatiles							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2-Hexanone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
2-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
4-Methyl-2-pentanone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C

Project ID: 16D3253 Phoenix I.D.: CF51978 Client ID: MW-104

RL/ Parameter Result **PQL** Units Dilution Date/Time Βv Reference ND 25 ug/L 1 03/17/20 МН SW8260C Acetone 1 ND 1.0 ug/L 03/17/20 MH SW8260C Acrylonitrile Benzene ND 0.70 ug/L 1 03/17/20 MH SW8260C ND 1.0 ug/L 1 03/17/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 03/17/20 MH SW8260C Bromochloromethane 1 Bromodichloromethane ND 0.50 ug/L 03/17/20 MH SW8260C 1 SW8260C ND 1.0 ug/L 03/17/20 **Bromoform** MH ND 1 Bromomethane 1.0 ug/L 03/17/20 MH SW8260C ug/L ND 5.0 1 03/17/20 MH SW8260C Carbon Disulfide 1 SW8260C Carbon tetrachloride ND 1.0 ug/L 03/17/20 MH Chlorobenzene ND 1.0 ug/L 1 03/17/20 MH SW8260C ND Chloroethane 1.0 ug/L 1 03/17/20 MH SW8260C ND 1.0 1 Chloroform ug/L 03/17/20 MH SW8260C ND 1 1.0 ug/L 03/17/20 MH SW8260C Chloromethane ND 1.0 ug/L 1 03/17/20 MH SW8260C cis-1,2-Dichloroethene cis-1,3-Dichloropropene ND 0.40 ug/L 1 03/17/20 MH SW8260C ND 0.50 ug/L 1 03/17/20 MH SW8260C Dibromochloromethane Dibromomethane ND 1.0 ug/L 1 03/17/20 MH SW8260C ND 1.0 ug/L 1 03/17/20 MH SW8260C Dichlorodifluoromethane ND 1 SW8260C 1.0 ug/L 03/17/20 MH Ethylbenzene ND 0.40 ug/L 1 03/17/20 MH SW8260C Hexachlorobutadiene Isopropylbenzene ND 1.0 ug/L 1 03/17/20 MH SW8260C m&p-Xylene ND 1.0 ug/L 1 03/17/20 МН SW8260C SW8260C Methyl ethyl ketone ND 5.0 ug/L 1 03/17/20 MH ND 1.0 ug/L 1 03/17/20 МН SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 03/17/20 МН SW8260C Methylene chloride ND 1.0 ug/L 1 03/17/20 МН SW8260C Naphthalene n-Butylbenzene ND 1.0 ug/L 1 03/17/20 MH SW8260C ND 1.0 ug/L 1 03/17/20 МН SW8260C n-Propylbenzene ND 1 SW8260C o-Xylene 1.0 ug/L 03/17/20 MH ND 1 SW8260C p-Isopropyltoluene 1.0 ug/L 03/17/20 MH ND 1 03/17/20 SW8260C sec-Butylbenzene 1.0 ug/L МН Styrene ND 1.0 ug/L 1 03/17/20 МН SW8260C ND 1 03/17/20 SW8260C tert-Butylbenzene 1.0 ug/L MH Tetrachloroethene ND 1.0 ug/L 1 03/17/20 МН SW8260C ND 2.5 SW8260C Tetrahydrofuran (THF) ug/L 1 03/17/20 MH ND 1.0 1 03/17/20 МН SW8260C ug/L Toluene ND 1 SW8260C **Total Xylenes** 1.0 ug/L 03/17/20 МН ND 1.0 ug/L 1 03/17/20 МН SW8260C trans-1,2-Dichloroethene 0.40 1 03/17/20 SW8260C trans-1,3-Dichloropropene ND ug/L MH ND 5.0 ug/L 1 03/17/20 МН SW8260C trans-1,4-dichloro-2-butene ND 1.0 ug/L 1 03/17/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 03/17/20 MH SW8260C Trichlorofluoromethane ND 1 0 ug/L 1 03/17/20 МН SW8260C Trichlorotrifluoroethane ug/L SW8260C ND 1.0 1 03/17/20 МН Vinyl chloride **QA/QC Surrogates** 95 % 1 03/17/20 МН 70 - 130 % % 1,2-dichlorobenzene-d4 92 % 1 03/17/20 MH 70 - 130 % % Bromofluorobenzene 03/17/20 70 - 130 % % Dibromofluoromethane 87 % 1 MH

Client ID: MW-104

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
% Toluene-d8	93		%	1	03/17/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	03/17/20	МН	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	МН	SW8260C (OXY)

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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:

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

Analysis Report

March 24, 2020

FOR: Attn: Matthew Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Matrix: WATER Collected by: 03/17/20

Location Code: BLCOMPANIES Received by: CP 03/17/20 16:16

Rush Request: 72 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data SDG ID: GCF51973

Phoenix ID: CF51979

Project ID: 16D3253 Client ID: TRIP BLANK

RL/

Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Volatiles							_
	ND	4.0	//	4	02/47/00	N 41 1	CMOCCOC
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,1-Dichloropropene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	03/17/20	МН	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
2-Hexanone	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
2-lisopropyltoluene	ND	1.0	ug/L	' 1	03/17/20	MH	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	5.0	=	1	03/17/20	MH	SW8260C
4-Methyl-2-pentanone	טאו	5.0	ug/L	ı	03/11/20	IVII ⁻¹	3002000

Project ID: 16D3253 Phoenix I.D.: CF51979
Client ID: TRIP BLANK

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Acetone	ND	25	ug/L	1	03/17/20	МН	SW8260C
Acrylonitrile	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Benzene	ND	0.70	ug/L	1	03/17/20	MH	SW8260C
Bromobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromochloromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	03/17/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Bromomethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Chloroform	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Chloromethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
cis-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	МН	SW8260C
Dibromochloromethane	ND	0.50	ug/L	1	03/17/20	МН	SW8260C
Dibromomethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Dichlorodifluoromethane	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Ethylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	03/17/20	МН	SW8260C
Isopropylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
m&p-Xylene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	03/17/20	МН	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Methylene chloride	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Naphthalene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
n-Propylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
o-Xylene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
p-Isopropyltoluene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
sec-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Styrene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Tetrachloroethene	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	03/17/20	МН	SW8260C
Toluene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Total Xylenes	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	03/17/20	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	03/17/20	MH	SW8260C
Trichloroethene	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
Trichlorotrifluoroethane	ND	1.0	ug/L	1	03/17/20	MH	SW8260C
	ND	1.0	ug/L	1	03/17/20	МН	SW8260C
Vinyl chloride	טאו	1.0	ug/L	ı	00/11/20	IVI□	3770200C
QA/QC Surrogates	97		%	1	03/17/20	M	70 - 130 %
% 1,2-dichlorobenzene-d4						MH	
% Bromofluorobenzene	93		%	1	03/17/20	MH	70 - 130 %
% Dibromofluoromethane	96		%	1	03/17/20	МН	70 - 130 %

Client ID: TRIP BLANK

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
i arameter	Result	I QL	Offits	Dilation	Date/Time	Dy	Reference
% Toluene-d8	93		%	1	03/17/20	MH	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	03/17/20	МН	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	03/17/20	MH	SW8260C (OXY)

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

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:

TRIP BLANK INCLUDED.

8260 Analysis:

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

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

Phyllis Shiller, Laboratory Director

March 24, 2020

Reviewed and Released by: Rashmi Makol, Project Manager



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

QA/QC Report

March 24, 2020

QA/QC Data

SDG I.D.: GCF51973

Parameter	Blank	Blk RL	Sample Result	Dup Result	Dup RPD	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
QA/QC Batch 522893 (mg/L), C	C Sam	ole No: (CF51552	(CF5197	4, CF5	1975,	CF5197	7)					
Mercury - Water Comment:	BRL	0.0002	<0.0002	<0.0002	NC	114			110			75 - 125	30
Additional Mercury criteria: LCS acceptance range for waters is 80-120% and for soils is 75-125%													
QA/QC Batch 522830 (mg/L), QC Sample No: CF50556 (CF51974, CF51975, CF51977)													
ICP Metals - Aqueous													
Antimony	BRL	0.005	< 0.005	< 0.005	NC	110	110	0.0	111			80 - 120	20
Arsenic	BRL	0.004	< 0.004	< 0.004	NC	104	104	0.0	106			80 - 120	20
Barium	BRL	0.002	0.032	0.033	3.10	104	103	1.0	104			80 - 120	20
Beryllium	BRL	0.001	< 0.001	< 0.001	NC	108	107	0.9	108			80 - 120	20
Cadmium	BRL	0.001	< 0.001	< 0.001	NC	108	107	0.9	107			80 - 120	20
Chromium	BRL	0.001	< 0.001	< 0.001	NC	104	103	1.0	104			80 - 120	20
Lead	BRL	0.002	< 0.002	< 0.002	NC	101	101	0.0	101			80 - 120	20
Nickel	BRL	0.001	< 0.001	< 0.001	NC	104	103	1.0	104			80 - 120	20
Selenium	BRL	0.010	< 0.010	< 0.010	NC	101	100	1.0	101			80 - 120	20
Silver	BRL	0.001	< 0.001	< 0.001	NC	99.9	98.8	1.1	100			80 - 120	20
Vanadium	BRL	0.002	< 0.002	< 0.002	NC	102	101	1.0	103			80 - 120	20
Zinc	BRL	0.004	< 0.004	< 0.004	NC	104	103	1.0	104			80 - 120	20
Comment:													
Additional: LCS acceptance range	is 80-12	0% MS a	cceptance	e range 7	5-125%								
QA/QC Batch 522992 (mg/L), C	C Sam	ole No: (CF52453	5X (CF5	1974, (CF5197	75, CF51	1977)					
ICP MS Metals - Aqueous	· }												
Thallium	BRL	0.0005	<0.0005	<0.0005	NC	103	104	1.0	104			80 - 120	20
Comment:													
Additional: LCS acceptance range is 80-120% MS acceptance range 75-125%.													



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QA/QC Report

QA/QC Data

March 24, 2020		QA/QC Data						SDG I.D.: GCF51973							
Parameter	Blank	Blk RL		LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits				
QA/QC Batch 522886 (ug/L), QC	Samp	e No: CF5197	4 (CF51974, CF5	51975, 0	CF51977	")									
MAEPH - Water				•		,									
C9-C18 Aliphatic Hydrocarbons 1*	ND	100		62	63	1.6				40 - 140	25				
C19-C36 Aliphatic Hydrocarbons 1*	ND	100		79	82	3.7				40 - 140	25				
C11-C22 Aromatic Hydrocarbons 1	ND	100		60	71	16.8				40 - 140	25				
C11-C22 Aromatic Hydrocarbons U	ND	100								40 - 140	25				
Total TPH 1,2*	ND	100		71	74	4.1				40 - 140	25				
C9 - Nonane	ND	10		38	36	5.4				40 - 140	25	1			
C-10 Decane	ND	10		50	48	4.1				40 - 140	25				
C12 - Dodecane	ND	10		56	55	1.8				40 - 140	25				
C14 - Tetradecane	ND	10		64	67	4.6				40 - 140	25				
C16 - Hexadecane	ND	10		74	78	5.3				40 - 140	25				
C18 - Octadecane	11	10		89	93	4.4				40 - 140	25				
C19 - Nonadecane	ND	10		83	86	3.6				40 - 140	25				
C20 - Eicosane	ND	10		84	87	3.5				40 - 140	25				
C22 - Docosane	ND	10		85	88	3.5				40 - 140	25				
C24 - Tetracosane	ND	10		86	90	4.5				40 - 140	25				
C26 - Hexacosane	ND	10		85	89	4.6				40 - 140	25				
C28 - Octacosane	ND	10		84	88	4.7				40 - 140	25				
C30 - Tricotane	ND	10		82	85	3.6				40 - 140	25				
C36 - Hexatriacontane	ND	10		41	45	9.3				40 - 140	25				
% 1-chlorooctadecane (aliphatic)	56	%		74	78	5.3				40 - 140	25				
% o-terphenyl (aromatic)	85	%		73	89	19.8				40 - 140	25				
% 2-Fluorobiphenyl (Fractionation)	110	%		90	104	14.4				40 - 140	25				
% 2-Bromonaphthalene (Fractionati	101	%		87	102	15.9				40 - 140	25				
% 2-Methylnaphthalene BT		%		0	0	NC				0 - 5					
% Naphthalene BT		%		0	0	NC				0 - 5					
Comment:															
Additional EPH fractionation criteria	a: Breakt	hrough criteria (l	BT) is 0 to 5%												
QA/QC Batch 522813 (ug/L), QC	Samp	e No: CF5101	2 (CF51974, CF5	1975, C	CF51977	')									
Semivolatiles - Water															
1,2,4,5-Tetrachlorobenzene	ND	3.5		75	80	6.5				40 - 140	20				
1,2,4-Trichlorobenzene	ND	3.5		74	83	11.5				40 - 140	20				
1,2-Dichlorobenzene	ND	1.0		74	79	6.5				40 - 140	20				
1,2-Diphenylhydrazine	ND	1.6		93	96	3.2				40 - 140	20				
1,3-Dichlorobenzene	ND	1.0		74	80	7.8				40 - 140	20				
1,4-Dichlorobenzene	ND	1.0		74	79	6.5				40 - 140	20				
2,4,5-Trichlorophenol	ND	1.0		92	98	6.3				30 - 130	20				
2,4,6-Trichlorophenol	ND	1.0		91	98	7.4				30 - 130	20				
2,4-Dichlorophenol	ND	1.0		85	91	6.8				30 - 130	20				
2,4-Dimethylphenol	ND	1.0		83	88	5.8				30 - 130	20				
2,4-Dinitrophenol	ND	1.0		92	102	10.3				30 - 130	20				

SDG I.D.: GCF51973

% % Blk **LCSD** LCS **RPD** LCS MS **MSD** MS Rec Blank RL **RPD** % % RPD Limits Limits % % Parameter 2,4-Dinitrotoluene ND 3.5 96 105 9.0 40 - 140 20 2,6-Dinitrotoluene ND 3.5 99 101 2.0 40 - 140 20 2-Chloronaphthalene ND 3.5 82 88 7.1 40 - 140 20 2-Chlorophenol ND 1.0 78 87 10.9 30 - 130 20 2-Methylphenol (o-cresol) ND 1.0 78 89 13.2 30 - 130 20 2-Nitroaniline ND 3.5 133 147 10.0 40 - 140 20 ND 1.0 83 102 20.5 30 - 130 20 2-Nitrophenol ND 1.0 97 30 - 130 3&4-Methylphenol (m&p-cresol) 83 15.6 20 3,3'-Dichlorobenzidine ND 48 40 - 140 5.0 44 8.7 20 3-Nitroaniline ND 5.0 102 105 2.9 40 - 140 20 4,6-Dinitro-2-methylphenol ND 1.0 92 101 9.3 30 - 130 20 ND 91 96 40 - 140 4-Bromophenyl phenyl ether 3.5 5.3 20 4-Chloro-3-methylphenol ND 1.0 94 102 8.2 30 - 130 20 4-Chloroaniline ND 3.5 78 82 5.0 40 - 140 20 4-Chlorophenyl phenyl ether ND 1.0 86 88 2.3 40 - 140 20 4-Nitroaniline ND 5.0 94 103 9.1 40 - 140 20 ND 4-Nitrophenol 1.0 100 112 11.3 30 - 130 20 ND 3.5 77 84 Acetophenone 8.7 40 - 140 20 ND 94 Aniline 3.5 84 11.2 40 - 140 20 Benzidine ND 4.5 92 104 12.2 40 - 140 20 ND 10 57 57 Benzoic acid 0.0 30 - 130 20 Benzyl butyl phthalate ND 1.5 94 104 10.1 40 - 140 20 ND 3.5 87 Bis(2-chloroethoxy)methane 82 5.9 40 - 140 20 Bis(2-chloroethyl)ether ND 1.0 72 81 11.8 40 - 140 20 Bis(2-chloroisopropyl)ether ND 1.0 72 80 10.5 40 - 140 20 ND 104 Bis(2-ethylhexyl)phthalate 1.5 96 8.0 40 - 140 20 Carbazole ND 5.0 91 96 5.3 40 - 140 20 ND 3.5 85 29 Dibenzofuran 4.6 40 - 140 20 ND 1.5 95 100 Diethyl phthalate 5.1 40 - 140 20 Dimethylphthalate ND 1.5 90 96 6.5 40 - 140 20 Di-n-butylphthalate ND 1.5 93 99 6.3 40 - 140 20 Di-n-octylphthalate ND 107 1.5 100 6.8 40 - 140 20 Hexachloroethane ND 3.5 75 83 10.1 40 - 140 20 ND Isophorone 3.5 83 86 3.6 40 - 140 20 N-Nitrosodi-n-propylamine ND 3.5 83 93 11.4 40 - 140 20 N-Nitrosodiphenylamine ND 3.5 81 84 3.6 40 - 140 20 ND 94 Pentachloronitrobenzene 5.0 100 6.2 40 - 140 20 Phenol ND 1.0 77 82 6.3 30 - 130 20 68 % 96 108 % 2,4,6-Tribromophenol 11.8 15 - 110 20 % 2-Fluorobiphenyl 75 % 75 77 2.6 30 - 130 20 62 78 % 2-Fluorophenol % 66 16.7 15 - 110 20 65 % Nitrobenzene-d5 % 76 82 7.6 30 - 130 20 65 % 71 83 15.6 % Phenol-d5 15 - 110 20 % Terphenyl-d14 75 % 81 84 3.6 30 - 130 20 Comment:

Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)

QA/QC Batch 522813 (ug/L), QC Sample No: CF51012 (CF51974, CF51975, CF51977)

Semivolatiles (SIM) - Water

2-Methylnaphthalene	ND	0.50	84	85	1.2	40 - 140	20
Acenaphthene	ND	0.50	108	109	0.9	40 - 140	20
Acenaphthylene	ND	0.50	103	106	2.9	40 - 140	20

SDG I.D.: GCF51973

Parameter	Blank	BIk RL	LC %		CSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
Anthracene	ND	0.50	10	2 1	105	2.9				40 - 140	20	
Benz(a)anthracene	ND	0.50	10	5 1	105	0.0				40 - 140	20	
Benzo(a)pyrene	ND	0.50	98	3 1	100	2.0				40 - 140	20	
Benzo(b)fluoranthene	ND	0.50	10	2 1	106	3.8				40 - 140	20	
Benzo(ghi)perylene	ND	0.50	10	8 1	110	1.8				40 - 140	20	
Benzo(k)fluoranthene	ND	0.50	98	3	99	1.0				40 - 140	20	
Chrysene	ND	0.50	10	5 1	107	1.9				40 - 140	20	
Dibenz(a,h)anthracene	ND	0.50	10	9 1	111	1.8				40 - 140	20	
Fluoranthene	ND	0.50	10	1 1	104	2.9				40 - 140	20	
Fluorene	ND	0.50	10	6 1	109	2.8				40 - 140	20	
Hexachlorobenzene	ND	0.50	90)	95	5.4				40 - 140	20	
Hexachlorobutadiene	ND	0.50	72	2	76	5.4				40 - 140	20	
Hexachlorocyclopentadiene	ND	0.50	30)	41	5.0				40 - 140	20	1
Indeno(1,2,3-cd)pyrene	ND	0.50	10	0 1	101	1.0				40 - 140	20	
Naphthalene	ND	0.50	78	3	79	1.3				40 - 140	20	
Nitrobenzene	ND	0.50	9		98	7.4				40 - 140	20	
N-Nitrosodimethylamine	ND	0.05	98	3 1	106	7.8				40 - 140	20	
Pentachlorophenol	ND	0.50	6		55	10.3				40 - 140	20	
Phenanthrene	ND	0.50	86)	91	2.2				40 - 140	20	
Pyrene	ND	0.50	10	8 1	112	3.6				40 - 140	20	
Pyridine	ND	0.50	7:	3	91	22.0				40 - 140	20	r
% 2,4,6-Tribromophenol	67	%	98	3	96	2.1				15 - 110	20	
% 2-Fluorobiphenyl	70	%	86)	89	0.0				40 - 140	20	
% 2-Fluorophenol	65	%	78	3	81	3.8				15 - 110	20	
% Nitrobenzene-d5	64	%	90)	88	2.2				40 - 140	20	
% Phenol-d5	71	%	9		94	3.2				15 - 110	20	
% Terphenyl-d14	72	%	90)	92	2.2				40 - 140	20	
Comment:												

Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)

QA/QC Batch 522853 (ug/L), QC Sample No: CF52086 (CF51974, CF51975, CF51977)

Semivolatiles by SIM, PAH - Water

Semivolatiles by Silvi,	PAH - Wa	<u>itei</u>							
2-Methylnaphthalene	ND	0.50	66	29	77.9	4	10 - 140	20	I,r
Acenaphthene	ND	0.50	170	72	81.0	4	10 - 140	20	l,r
Acenaphthylene	ND	0.10	166	71	80.2	4	10 - 140	20	l,r
Anthracene	ND	0.10	158	67	80.9	4	10 - 140	20	l,r
Benz(a)anthracene	ND	0.02	175	73	82.3	4	10 - 140	20	l,r
Benzo(a)pyrene	ND	0.02	163	69	81.0	4	10 - 140	20	l,r
Benzo(b)fluoranthene	ND	0.02	159	69	78.9	4	10 - 140	20	l,r
Benzo(ghi)perylene	ND	0.02	185	81	78.2	4	10 - 140	20	l,r
Benzo(k)fluoranthene	ND	0.02	149	63	81.1	4	10 - 140	20	l,r
Chrysene	ND	0.02	165	70	80.9	4	10 - 140	20	l,r
Dibenz(a,h)anthracene	ND	0.02	171	73	80.3	4	10 - 140	20	l,r
Fluoranthene	ND	0.50	155	64	83.1	4	10 - 140	20	l,r
Fluorene	ND	0.10	165	72	78.5	4	10 - 140	20	l,r
Indeno(1,2,3-cd)pyrene	ND	0.02	168	71	81.2	4	10 - 140	20	l,r
Naphthalene	ND	0.50	94	41	78.5	4	10 - 140	20	r
Phenanthrene	ND	0.50	151	63	82.2	4	10 - 140	20	I,r
Pyrene	ND	0.07	163	67	83.5	4	10 - 140	20	I,r
% 2-Fluorobiphenyl	77	%	72	77	6.7	4	10 - 140	20	
% Nitrobenzene-d5	75	%	63	70	10.5	4	10 - 140	20	
% Terphenyl-d14	67	%	65	65	0.0	4	10 - 140	20	

SDG I.D.: GCF51973

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
Comment:											
Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)											
QA/QC Batch 522985 (ug/L), C	C Samp	le No: CF51399 (CF51973, CF	51974. (CF51975	5. CF51	976. C	F51977	. CF51	978, CF	51979)	
Volatiles - Water		(0. 0)	, .		.,	, -		,	,	,	
1,1,1,2-Tetrachloroethane	ND	1.0	92	96	4.3	94	94	0.0	70 120	20	
1,1,1,2-retrachioroethane	ND ND	1.0	92 86	96 89	4.3 3.4	94 87	94 87	0.0	70 - 130 70 - 130	30 30	
1,1,2,2-Tetrachloroethane	ND ND	0.50	96	09 107	3.4 10.8	101	106	4.8	70 - 130	30	
1,1,2-Trichloroethane	ND	1.0	88	98	10.8	95	97	2.1	70 - 130	30	
1,1-Dichloroethane	ND	1.0	91	92	1.1	93 87	89	2.3	70 - 130	30	
1,1-Dichloroethene	ND	1.0	93	97	4.2	91	94	3.2	70 - 130	30	
1,1-Dichloropropene	ND	1.0	83	86	3.6	89	90	1.1	70 - 130	30	
1,2,3-Trichlorobenzene	ND	1.0	92	101	9.3	86	100	15.1	70 - 130	30	
1,2,3-Trichloropropane	ND	1.0	80	88	9.5	85	89	4.6	70 - 130	30	
1,2,4-Trichlorobenzene	ND	1.0	91	100	9.4	89	99	10.6	70 - 130	30	
1,2,4-Trimethylbenzene	ND	1.0	97	98	1.0	97	98	1.0	70 - 130	30	
1,2-Dibromo-3-chloropropane	ND	1.0	95	107	11.9	91	98	7.4	70 - 130	30	
1,2-Dibromoethane	ND	1.0	88	95	7.7	88	91	3.4	70 - 130	30	
1,2-Dichlorobenzene	ND	1.0	93	99	6.3	95	98	3.1	70 - 130	30	
1,2-Dichloroethane	ND	1.0	84	92	9.1	87	89	2.3	70 - 130	30	
1,2-Dichloropropane	ND	1.0	93	99	6.3	98	99	1.0	70 - 130	30	
1,3,5-Trimethylbenzene	ND	1.0	97	98	1.0	98	99	1.0	70 - 130	30	
1,3-Dichlorobenzene	ND	1.0	94	98	4.2	95	96	1.0	70 - 130	30	
1,3-Dichloropropane	ND	1.0	88	95	7.7	91	94	3.2	70 - 130	30	
1,4-Dichlorobenzene	ND	1.0	92	97	5.3	94	96	2.1	70 - 130	30	
1,4-dioxane	ND	100	89	90	1.1	69	71	2.9	40 - 160	30	
2,2-Dichloropropane	ND	1.0	87	87	0.0	79	79	0.0	70 - 130	30	
2-Chlorotoluene	ND	1.0	99	102	3.0	101	101	0.0	70 - 130	30	
2-Hexanone	ND	5.0	77	92	17.8	76	85	11.2	40 - 160	30	
2-Isopropyltoluene	ND	1.0	105	105	0.0	106	106	0.0	70 - 130	30	
4-Chlorotoluene	ND	1.0	96	97	1.0	97	98	1.0	70 - 130	30	
4-Methyl-2-pentanone	ND	5.0	81	94	14.9	88	93	5.5	40 - 160	30	
Acetone	ND	5.0	83	86	3.6	81	89	9.4	40 - 160	30	
Acrylonitrile	ND	5.0	85	98	14.2	85	92	7.9	70 - 130	30	
Benzene	ND	0.70	91	94	3.2	111	110	0.9	70 - 130	30	
Bromobenzene	ND	1.0	94	100	6.2	95	97	2.1	70 - 130	30	
Bromochloromethane	ND	1.0	85	92	7.9	84	85	1.2	70 - 130	30	
Bromodichloromethane	ND	0.50	91	98	7.4	96	97	1.0	70 - 130	30	
Bromoform	ND	1.0	92	105	13.2	94	99	5.2	70 - 130	30	
Bromomethane	ND	1.0	106	109	2.8	74	93	22.8	40 - 160	30	
Carbon Disulfide	ND	1.0	95	94	1.1	89	93	4.4	70 - 130	30	
Carbon tetrachloride	ND	1.0	100	102	2.0	104	103	1.0	70 - 130	30	
Chlorobenzene	ND	1.0	93	97	4.2	97	97	0.0	70 - 130	30	
Chloroethane	ND	1.0	91	88	3.4	80	84	4.9	70 - 130	30	
Chloroform	ND	1.0	91	86	5.6	89	83	7.0	70 - 130	30	
Chloromethane	ND	1.0	90	92	2.2	75	80	6.5	40 - 160	30	
cis-1,2-Dichloroethene	ND	1.0	89	92	3.3	85	89	4.6	70 - 130	30	
cis-1,3-Dichloropropene	ND	0.40	91	100	9.4	95	99	4.1	70 - 130	30	
Dibromochloromethane	ND	0.50	95	104	9.0	97	99	2.0	70 - 130	30	
Dibromomethane	ND	1.0	88	94	6.6	92	95	3.2	70 - 130	30	
Dichlorodifluoromethane	ND	1.0	81	81	0.0	63	64	1.6	40 - 160	30	
Ethyl ether	ND	1.0	93	103	10.2	92	96	4.3	70 - 130	30	

OA	Ω	Data
$\nabla \alpha$	\sim	Data

SDG I.D.: GCF51973

Ethylbenzene	Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
Hexaphorbulatedine ND 0.40 100 846 97 101 20 20 101 20 70 101 20 20	Ethylbenzene	ND	1.0	98	99	1.0	101	101	0.0	70 - 130	30	
Methylethylethylethone Ni		ND	0.40	109	100	8.6	99	101	2.0	70 - 130	30	
Methyle thely ketone	Isopropylbenzene	ND	1.0	98	100	2.0	102	102	0.0	70 - 130	30	
Methyle thay katone ND	· · · · ·	ND		95	97	2.1	98	99		70 - 130	30	
Methylen chloride (MTBE)	· · ·	ND		77	90	15.6	74	88	17.3	40 - 160	30	
Methylakene chloride		ND		84	94	11.2	79			70 - 130	30	
Namipulaene N.D 1.0 1.				88	91	3.4	82	86		70 - 130	30	
n-Brogylbanzene ND 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0	_	ND		90	104	14.4	86	104		70 - 130	30	
n-Progrighenzene	•	ND	1.0	104	105	1.0	106	108	1.9	70 - 130	30	
Section No 1.0 1		ND	1.0	98	99	1.0	100	102	2.0	70 - 130	30	
P-Isophyghlouene	: -	ND	1.0	98	101	3.0	101	101	0.0	70 - 130	30	
Second No. 1.0 1	-	ND	1.0	101	100	1.0	102	103	1.0	70 - 130	30	
Styrene		ND	1.0	105	106	0.9	110	110	0.0	70 - 130	30	
Refi-Bulylbenzene N.D. 1.0 1		ND	1.0	95	99	4.1	98	100	2.0	70 - 130	30	
Petrachloroethene	=	ND	1.0	98	99	1.0	101	101	0.0	70 - 130	30	
Tolluene ND 1.0 98 102 4.0 105 106 0.9 70 - 130 30 1 1 1 1 1 1 1 1 1	=	ND	1.0	87	93	6.7	96	97	1.0	70 - 130	30	
trans-1,2-Dichloroethene ND 1.0 92 94 2.2 88 92 4.4 70-130 30 trans-1,3-Dichloropropene ND 0.40 87 97 10.9 90 94 4.3 70-130 30 trans-1,3-Dichloropropene ND 5.0 149 165 10.2 10 98 92 31.2 160 9.8 70-130 30 Image: No 10 10 99 92 32.3 95 90 0.0 70-130 30 71-710 70-710 99 100 1.0 98 99 100 1.0 98 91 90 90 90 100 9	Tetrahydrofuran (THF)	ND	2.5	78	95	19.7	76	84	10.0	70 - 130	30	
trans-1,3-Dichloropropene ND 0.40 87 97 10.9 90 43 70-130 30 Intrans-1,4-dichloro-2-butene ND 5.0 148 165 10.2 148 160 98 70-130 30 Image: North Introduction (Introduction (Introduct	Toluene	ND	1.0	98	102	4.0	105	106	0.9	70 - 130	30	
trans-1,4-dichloro-2-butene ND 5.0 149 165 10.2 145 160 9.8 70-130 30 Image: Not of the properties of the pr	trans-1,2-Dichloroethene	ND	1.0	92	94	2.2	88	92	4.4	70 - 130	30	
Trichloroethene ND 1.0 89 92 3.3 95 95 0.0 70-130 30 Trichlorofurormethane ND 1.0 92 94 2.2 89 91 2.2 70-130 30 Trichlorotifluroethane ND 1.0 101 105 3.9 95 60 70-130 30 % Invigority chioride ND 1.0 99 100 102 2.0 102 102 0.0 70-130 30 % Bromofluorobenzene-d4 95 % 100 102 2.0 102 102 102 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 30 70-130 </td <td>trans-1,3-Dichloropropene</td> <td>ND</td> <td>0.40</td> <td>87</td> <td>97</td> <td>10.9</td> <td>90</td> <td>94</td> <td>4.3</td> <td>70 - 130</td> <td>30</td> <td></td>	trans-1,3-Dichloropropene	ND	0.40	87	97	10.9	90	94	4.3	70 - 130	30	
Trichlorofluoromethane ND 1.0 92 94 2.2 89 91 2.2 70-130 30 Trichlorotifluoroethane ND 1.0 101 105 3.9 94 96 2.1 70-130 30 Vinyl chloride ND 1.0 99 100 1.0 86 91 5.0 70-130 30 % 1.2-dichlorobenzene-d4 95 % 98 99 1.0 98 101 3.0 70-130 30 % Dibromofluoromethane 92 % 90 93 3.3 83 87 4.7 70-130 30 % Dibromofluoromethane 93 % 100 100 102 102 102 70-130 30 A/L/CC Batch \$23354 (ug/L/), CC Sampler No: CF\$52176 (CF\$51975 (20X) V V 70-130 30 V Volatiles - Water Trichlorotifluoroethane ND 1.0 62 83 29.0 V		ND	5.0	149	165	10.2	145	160	9.8	70 - 130	30	l,m
Trichlorotrifluoroethane ND 1.0 101 105 3.9 94 96 2.1 70-130 30 Vinyl chloride ND 1.0 99 100 1.0 86 91 5.6 70-130 30 % 1,2-dichlorobenzene-d4 95 % 100 102 2.0 102 102 0.0 70-130 30 % Bromofluorobenzene 91 % 98 99 1.0 98 101 3.0 70-130 30 % Bromofluorobenzene 92 % 90 93 3.3 83 87 4.7 70-130 30 % Toluene-d8 93 % 100 100 100 102 102 102 70-130 30 Comment: Additional 8260 criteria: 10% of components and an	Trichloroethene	ND	1.0	89	92	3.3	95	95	0.0	70 - 130	30	
Niny chloride	Trichlorofluoromethane	ND	1.0	92	94	2.2	89	91	2.2	70 - 130	30	
% 1,2-dichlorobenzene-d4 95 % 100 102 2.0 102 102 102 102 102 0.0 70 - 130 30 % Bromofluorobenzene 91 % 98 99 1.0 98 101 3.0 70 - 130 30 % Dibromofluorobenzene 93 % 100 100 100 100 102 102 102 0.0 70 - 130 30 Comment: Additional 8260 criteria: 10% of compounds carb be outside of acceptance criteria as long as recovery is 10%. CA/QC Batch 523354 (ug/L), QC Sample No: CF52176 (CF51975 (20X)) Volatiles - Water Trichlorotifiluoroethane ND 1.0 62 83 29.0 - 70 - 130 30 - % 12-dichlorobenzene-d4 102 % 100 101 1.0 - 70 - 130 30 - % Dibromofluoromethane 103 % 99 99 0.0 - 10 70 - 130 30 </td <td>Trichlorotrifluoroethane</td> <td>ND</td> <td>1.0</td> <td>101</td> <td>105</td> <td>3.9</td> <td>94</td> <td>96</td> <td>2.1</td> <td>70 - 130</td> <td>30</td> <td></td>	Trichlorotrifluoroethane	ND	1.0	101	105	3.9	94	96	2.1	70 - 130	30	
Se Formoriluorobenzene 91 % 98 99 1.0 98 101 3.0 70-130 30 % Dibromofluorobenzene 92 % 90 93 3.3 83 87 4.7 70-130 30 % Toluene-d8 93 % 100 100 100 100 102 102 0.0 70-130 30 70-130 30 Comment: Additional 8260 criteria: 10% of compounds: car be outside of acceptance criteria as lows: are recovery is 1000. 100 <td>Vinyl chloride</td> <td>ND</td> <td>1.0</td> <td>99</td> <td>100</td> <td>1.0</td> <td>86</td> <td>91</td> <td>5.6</td> <td>70 - 130</td> <td>30</td> <td></td>	Vinyl chloride	ND	1.0	99	100	1.0	86	91	5.6	70 - 130	30	
% Dibromofluoromethane 92 % 90 93 3.3 83 87 4.7 70-130 30 % Toluene-d8 93 % 100 100 0.0 102 102 0.0 70-130 30 Comment: Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. CAJA/QC Batch 523354 (ug/L), QC Sample No: CF52176 (CF51975 (20X)) Volatiles - Water Trichlorotrifluoroethane ND 1.0 62 83 29.0 2 2 70-130 30 1 % 1.2-dichlorobenzene-d4 102 % 100 101 1.0 2 70-130 30 1 % Bromofluoromethane 103 % 100 101 1.0 2 70-130 30 1 % Dibromofluoromethane 103 % 100 101 1.0 2 70-130 30 1 2 70-130 30 1 1 1	% 1,2-dichlorobenzene-d4	95	%	100	102	2.0	102	102	0.0	70 - 130	30	
% Toluene-d8 Comment: 93 % 100 100 100 102 102 102 70-130 30 Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as receivery is 10%. QA/QC Batch 523354 (ug/L), QC Sample No: CF52176 (CF51975 (20X)) Volatiles - Water Trichlorotrifluoroethane ND 1.0 62 83 29.0 1 70-130 30 1 % 1,2-dichlorobenzene-d4 102 % 100 101 1.0 1 70-130 30 1 % Bromofluorobenzene 96 % 99 99 99 90 0.0 70-130 30 % Dibromofluoromethane 103 % 100 101 1.0 70-130 30 % Dibromofluoromethane 103 % 100 101 1.0 70-130 30 % Toluene-d8 101 % 9 9	% Bromofluorobenzene	91	%	98	99	1.0	98	101	3.0	70 - 130	30	
Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. CAP/QC Batch 523354 (ug/L), QC Sample No: CF52176 (CF51975 (20X)) Volatiles - Water	% Dibromofluoromethane	92	%	90	93	3.3	83	87	4.7	70 - 130	30	
QA/QC Batch 523354 (ug/L), QC Sample No: CF52176 (CF51975 (20X)) Volatiles - Water Trichlorotrifluoroethane ND 1.0 62 83 29.0 70 - 130 30 1 % 1,2-dichlorobenzene-d4 102 % 100 101 1.0 70 - 130 30 1 % Bromofluorobenzene 96 % 99 99 0.0 70 - 130 30 1 <td></td> <td>93</td> <td>%</td> <td>100</td> <td>100</td> <td>0.0</td> <td>102</td> <td>102</td> <td>0.0</td> <td>70 - 130</td> <td>30</td> <td></td>		93	%	100	100	0.0	102	102	0.0	70 - 130	30	
Volatiles - Water Trichlorotrifluoroethane ND 1.0 62 83 29.0 70 - 130 30 1 % 1,2-dichlorobenzene-d4 102 % 100 101 1.0 70 - 130 30 1 % Bromofluorobenzene 96 % 99 99 0.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 <td>Additional 8260 criteria: 10% of o</td> <td>compounds</td> <td>can be outside of acceptance</td> <td>criteria as lor</td> <td>ng as reco</td> <td>overy is</td> <td>10%.</td> <td></td> <td></td> <td></td> <td></td> <td></td>	Additional 8260 criteria: 10% of o	compounds	can be outside of acceptance	criteria as lor	ng as reco	overy is	10%.					
Volatiles - Water Trichlorotrifluoroethane ND 1.0 62 83 29.0 70 - 130 30 1 % 1,2-dichlorobenzene-d4 102 % 100 101 1.0 70 - 130 30 1 % Bromofluorobenzene 96 % 99 99 0.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 70 - 130 30 1 1 1.0 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 1.0 1 <td>QA/QC Batch 523354 (ug/L), (</td> <td>QC Samp</td> <td>le No: CF52176 (CF51975 (</td> <td>(20X))</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	QA/QC Batch 523354 (ug/L), (QC Samp	le No: CF52176 (CF51975 ((20X))								
% 1,2-dichlorobenzene-d4 102 % 100 101 1.0 70 - 130 30 % Bromofluorobenzene 96 % 99 99 0.0 70 - 130 30 % Dibromofluoromethane 103 % 100 101 1.0 70 - 130 30 % Toluene-d8 101 % 99 99 0.0 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 compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichloroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30 <td>Volatiles - Water</td> <td>·</td> <td></td>	Volatiles - Water	·										
% Bromofluorobenzene 96 % 99 99 0.0 70 - 130 30 % Dibromofluoromethane 103 % 100 101 1.0 70 - 130 30 ** Toluene-d8 101 % 99 99 0.0 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 compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichloroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30												I
% Dibromofluoromethane 103 % 100 101 1.00 70 - 130 30 % Toluene-d8 101 % 99 99 0.0 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 compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichlorotrifluoroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30	·											
% Toluene-d8 Comment: 101 % 99 99 0.0 70 - 130 30 A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate. Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichlorotrifluoroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 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 compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene												
Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water Trichloroethene	Comment:					0.0				70 - 130	30	
QA/QC Batch 523171 (ug/L), QC Sample No: CF52205 (CF51974 (5X) , CF51975 (5X) , CF51977 (5X)) Volatiles - Water ND 1.0 93 91 2.2 70 - 130 30 Trichloroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30	•		•	•	•	m!-	100/					
Volatiles - Water Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichlorotrifluoroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30		•	•		•	,		X))				
Trichloroethene ND 1.0 93 91 2.2 70 - 130 30 Trichlorotrifluoroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30		ao oamp	01101010102200 (0101771)	(07,07,01,01)	,,,,	, 0, 0,	. , , , , (0	,,,,				
Trichlorotrifluoroethane ND 1.0 102 104 1.9 70 - 130 30 % 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30		NID	1.0	00	01	2.2				70 100	20	
% 1,2-dichlorobenzene-d4 96 % 102 101 1.0 70 - 130 30 % Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30												
% Bromofluorobenzene 92 % 98 99 1.0 70 - 130 30												
% Dibromofluoromethane 89 % 86 89 3.4 70 - 130 30 % Toluene-d8 92 % 102 101 1.0 70 - 130 30												

SDG I.D.: GCF51973

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

Comment:

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

Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%.

QA/QC Batch 523130 (ug/L), QC Sample No: CF51974 (CF51974, CF51975, CF51977)

Volatile Petroleum Hydrocarbons - Water

Unadjusted C5-C8 Aliphatics (*1)	ND	100	104	105	1.0	NC	NC	NC	70 - 130	20
Unadjusted C9-C12 Aliphatics (*1)	ND	100	95	97	2.1	NC	NC	NC	70 - 130	20
C5-C8 Aliphatic Hydrocarbons *1,2	ND	100	104	105	1.0	NC	NC	NC	70 - 130	20
C9-C12 Aliphatic Hydrocarbons *1,	ND	100	95	97	2.1	NC	NC	NC	70 - 130	20
C9-C10 Aromatic Hydrocarbons *1	ND	100	90	92	2.2	NC	NC	NC	70 - 130	20
Benzene	ND	1.0	92	92	0.0	95	95	0.0	70 - 130	20
Ethyl Benzene	ND	1.0	90	92	2.2	94	94	0.0	70 - 130	20
MTBE	ND	1.0	89	90	1.1	92	94	2.2	70 - 130	20
Naphthalene	ND	5.0	86	86	0.0	89	89	0.0	70 - 130	20
Toluene	ND	1.0	90	92	2.2	95	94	1.1	70 - 130	20
m,p-Xylenes	ND	2.0	91	92	1.1	95	96	1.0	70 - 130	20
o-Xylene	ND	1.0	89	91	2.2	92	93	1.1	70 - 130	20
% 2,5-Dibromotoluene (PID)	91	%	90	86	4.5	85	86	1.2	70 - 130	20
Comment:										

A blank MS/MSD was analyzed with this batch.

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

March 24, 2020

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

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

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

Tuesday, March 24, 2020

Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report GCF51973 - BLCOMPANIES

State: MA

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	Criteria	Units
CF51973	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51973	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51973	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51973	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51973	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51973	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51973	\$MCPADD-WM	I 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51973	\$MCPADD-WM	I 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51974	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51974	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51974	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51974	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51974	\$8260GWR	Trichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	20	1.0	5	5	ug/L
CF51974	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51974	\$8260GWR	Trichloroethene	MA / GROUNDWATER STANDARDS / GW-1	20	1.0	5	5	ug/L
CF51974	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51974	\$8270-SIMFSR	Benzoic acid	MA / CAM Protocol / SVOA AQ RL	ND	51		10	ug/L
CF51974	\$MCPADD-WM	I 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51974	\$MCPADD-WM	I 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51975	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51975	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51975	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51975	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51975	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51975	\$8260GWR	Trichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	44	5.0	5	5	ug/L
CF51975	\$8260GWR	Tetrachloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	7.0	1.0	5	5	ug/L
CF51975	\$8260GWR	Tetrachloroethene	MA / GROUNDWATER STANDARDS / GW-1	7.0	1.0	5	5	ug/L
CF51975	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51975	\$8260GWR	Trichloroethene	MA / GROUNDWATER STANDARDS / GW-1	44	5.0	5	5	ug/L
CF51975	\$8270-SIMFSR	Benzoic acid	MA / CAM Protocol / SVOA AQ RL	ND	51		10	ug/L
CF51975	\$8270-SIMR	Benzo(a)pyrene	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.21	0.2	0.2	ug/L
CF51975	\$8270-SIMR	Benzo(a)pyrene	MA / GROUNDWATER STANDARDS / GW-1	ND	0.21	0.2	0.2	ug/L
CF51975	\$MCPADD-WM	I 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51975	\$MCPADD-WM	I 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51976	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51976	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51976	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51976	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51976	\$8260GWR	Trichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	5.8	1.0	5	5	ug/L
CF51976	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L

RL

Analysis

Tuesday, March 24, 2020

Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report

State: MA

GCF51973 - BLCOMPANIES

SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	Criteria	Units
CF51976	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51976	\$8260GWR	Trichloroethene	MA / GROUNDWATER STANDARDS / GW-1	5.8	1.0	5	5	ug/L
CF51976	\$MCPADD-WM	1 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51976	\$MCPADD-WM	1 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51977	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51977	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51977	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51977	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51977	\$8260GWR	Trichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	34	5.0	5	5	ug/L
CF51977	\$8260GWR	Tetrachloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	5.4	1.0	5	5	ug/L
CF51977	\$8260GWR	cis-1,2-Dichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	30	1.0	20	20	ug/L
CF51977	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51977	\$8260GWR	Tetrachloroethene	MA / GROUNDWATER STANDARDS / GW-1	5.4	1.0	5	5	ug/L
CF51977	\$8260GWR	Trichloroethene	MA / GROUNDWATER STANDARDS / GW-1	34	5.0	5	5	ug/L
CF51977	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51977	\$8270-SIMFSR	Benzoic acid	MA / CAM Protocol / SVOA AQ RL	ND	55		10	ug/L
CF51977	\$8270-SIMFSR	3&4-Methylphenol (m&p-cresol)	MA / CAM Protocol / SVOA AQ RL	ND	11		10	ug/L
CF51977	\$MCPADD-WM	I 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51977	\$MCPADD-WM	1 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51978	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51978	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51978	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51978	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51978	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51978	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51978	\$MCPADD-WM	1 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51978	\$MCPADD-WM	1 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
CF51979	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF51979	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51979	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF51979	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF51979	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF51979	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF51979	\$MCPADD-WM	I 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF51979	\$MCPADD-WM	1 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	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 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.

RL

Analysis

		Mas	ssDE	P Analytic	cal P	rotocol Ce	ertific	ation Forn	1		
Labo	ratory Na	ame: Phoer	nix En	vironmental	Labora	atories, Inc.	Proje	ct #:			
Proje	ect Locat	ion: 16D32	253				RTN:				
This F	orm provid	des certification	ns for	the following o	data set	: [list Laborate	ory San	nple ID Number(s)]		
CF519	973, CF5197	74, CF51975, C	F5197	6, CF51977, CF	51978,	CF51979					
Matric		oundwater/Surfa			Sedimen	t Drinkir	ıg Wate	r 🗌 Air	✓ 0	ther: \	WATER
8260 V		7470/7471 Ho		MassDEP VPH	4	8081 Pesticide	9	7196 Hex Cr		MassD	EP APH
CAM II		CAM III B	✓	CAM IV A	· •	CAM V B		CAM VI B		CAM IX	
8270 S CAM II		7010 Metals CAM III C		MassDEP EPH CAM IV B	∀	8151 Herbicide CAM V C	es 🗌	8330 Explosives CAM VIII A		TO-15 CAM IX	
6010 M CAM II		6020 Metals CAM III D		8082 PCB CAM V A		9012 Total Cyanide/PAC CAM V1 A		6860 Perchlorat CAM VIII B			
	Affirmat	ive response	s to c	uestions A tl	nrough	F are require	ed for	"Presumptive	Certa	inty" s	tatus
Α	Chain-of-laboratory	Custody, prop /, and prepare	erly p ed/ana	a condition coreserved (inclinated in a condition of the	uding to	emperature*) olding times?	in the fi (* see r	ield or narrative)		Yes	✓ No
В		analytical me CAM protocol		s) and all asso owed?	ciated	QC requireme	ents sp	ecified in the	✓	Yes	\square No
С		CAM protocol		actions and ar plemented for					✓	Yes	□No
D	CAM VII A		suran	mply with all to be and Quality pata"?					✓	Yes	□No
Е	significan modificati	t modification ons).	(s)? (ods only: Wa refer to the ind only: Was the	dividual	method(s) fo	r a list o	of significant		Yes Yes	□ No
	method?		tillous	orny. Was the	Compi	cic analyte no	торог	ica for cacif		162	
F	conforma		d and	tocol QC and evaluated in a ough E)?				ling all "No"	✓	Yes	□ No
	Res	ponses to qu	estio	ns G, H and I	below	is required fo	or "Pre	sumptive Cert	tainty	' statu	S
G		reporting lim		r below all CA	М герс	orting limits sp	ecified	in the		Yes	☑ No
repres	entativene	ss requiremer	ts des	cribed in 310 C	MR 40.	1056(2)(k) and	d WSC-		data u	sability	and
Н	See Secti	ions: EPH, S\	/OA, S	andards specif SVOASIM, VO	A Narra	ations .				Yes	✓ No
I	Were resprotocol(s	s)?		complete ana		,				Yes	□ No
respor	nsible for o	d, attest unde	the part	ains and penal	ties of p	perjury that, ba	sed up	on my personal report is, to the	inqui		
							Date: 1	Гuesday, Marc	ch 24	2020	
	norized nature: -	Ros	hu	i nako	H	Printed Na	ame: F	Rashmi Makol			
gi						Pos	ition: F	Project Manag	er		



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MCP Certification Report

March 24, 2020 SDG I.D.: GCF51973

SDG Comments

8260 Analysis:

1,2-Dibromoethane doesn't meet GW-1 criteria, this compound is analyzed by GC/FID to achieve this criteria.

1,4-Dioxane doesn't meet GW-1 criteria, this compound is analyzed by 8270SIM to achieve this criteria.

Phoenix reporting levels may exceed those referenced in the CAM protocol. Please refer to criteria sheet for comparisons to requested MCP standards

Temperature above 6C:

The samples were received in a cooler with ice packs. No significant bias is suspected.

EPH Narration

Were all QA/QC performance criteria specified in the MADEP document CAM achieved? No.

QC Batch 522886 (Samples: CF51974, CF51975, CF51977): -----

A trace amount of an analyte was found in blank but were not reported in the sample(s), therefore no bias is suspected. (C18 - Octadecane)

One or more analytes is below the method criteria. A low bias for these analytes is possible. (C9 - Nonane)

Instrument:

AU-FID3 03/20/20-1

Adam Werner, Chemist 03/20/20

CF51975 (1X)

No significant modifications were made to the EPH method, as specified in Section 11.3 of the method.

The initial calibration (AR0102BI) RSD for the compound list was less than 20% except for the following compounds: None. The continuing calibration %D for the compound list was less than 25% except for the following compounds: None.

AU-FID4 03/17/20-1

Adam Werner, Chemist 03/17/20

CF51974 (1X), CF51977 (1X)

The initial calibration (AL0219BI) RSD for the compound list was less than 20% except for the following compounds: None. The continuing calibration %D for the compound list was less than 25% except for the following compounds: None.

AU-FID4 03/20/20-1

Adam Werner, Chemist 03/20/20

CF51975 (1X)

The initial calibration (AL0219BI) RSD for the compound list was less than 20% except for the following compounds: None. The continuing calibration %D for the compound list was less than 25% except for the following compounds: None.

QC (Batch Specific):

Batch 522886 (CF51974)

CF51974, CF51975, CF51977

All LCS recoveries were within 40 - 140 with the following exceptions: C9 - Nonane(38%) All LCSD recoveries were within 40 - 140 with the following exceptions: C9 - Nonane(36%)

All LCS/LCSD RPDs were less than 25% with the following exceptions: None. Additional EPH fractionation criteria: Breakthrough criteria (BT) is 0 to 5%

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MCP Certification Report

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EPH Narration

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

Mercury Narration

Were all QA/QC performance criteria specified in the analytical method achieved? Yes.

Instrument:

MERLIN 03/18/20 08:00

Rick Schweitzer, Chemist 03/18/20

CF51974, CF51975, CF51977

The method preparation blank, ICB, and CCBs contain all of the acids and reagents as the samples.

The initial calibration met all criteria including a standard run at or below the reporting level.

All calibration verification standards (ICV, CCV) met criteria.

All calibration blank verification standards (ICB, CCB) met criteria.

The matrix spike sample is used to identify spectral interference for each batch of samples, if within 85-115%, no interference is observed and no further action is taken.

The following Initial Calibration Verification (ICV) compounds did not meet criteria: None.

The following Continuing Calibration Verification (CCV) compounds did not meet criteria: None.

QC (Batch Specific):

Batch 522893 (CF51552)

CF51974, CF51975, CF51977

All LCS recoveries were within 75 - 125 with the following exceptions: None.

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

ICP Metals Narration

Were all QA/QC performance criteria specified in the analytical method achieved? Yes.

Instrument:

BLUE 03/18/20 08:26

Cindy Pearce, Chemist 03/18/20

CF51974, CF51975, CF51977

The initial calibration met criteria.

The continuing calibration standards met criteria for all the elements reported. The linear range is defined daily by the calibration range.

The continuing calibration blanks were less than the reporting level for the elements reported.

The ICSA and ICSAB were analyzed at the beginning and end of the run and were within criteria. The linear range is defined daily by the calibration range.

The following Initial Calibration Verification (ICV) compounds did not meet criteria: None.

The following Continuing Calibration Verification (CCV) compounds did not meet criteria: None.

The following ICP Interference Check (ICSAB) compounds did not meet criteria: None.

QC (Batch Specific):

Batch 522830 (CF50556)

CF51974, CF51975, CF51977

All LCS recoveries were within 80 - 120 with the following exceptions: None.



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Certification Report

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ICP Metals Narration

All LCSD recoveries were within 80 - 120 with the following exceptions: None. All LCS/LCSD RPDs were less than 20% with the following exceptions: None. Additional: LCS acceptance range is 80-120% MS acceptance range 75-125%.

ICPMS Metals Narration

Were all QA/QC performance criteria specified in the analytical method achieved? Yes.

Instrument:

ICPMS 03/19/20 14:53

Cindy Pearce, Chemist 03/19/20

CF51974, CF51975, CF51977

The linear range is defined daily by the calibration range.

The following Initial Calibration Verification (ICV) compounds did not meet criteria: None.

The following Continuing Calibration Verification (CCV) compounds did not meet criteria: None.

The following samples did not meet internal standard criteria: None.

QC (Batch Specific):

Batch 522992 (CF52453)

CF51974, CF51975, CF51977

All LCS recoveries were within 80 - 120 with the following exceptions: None.

All LCSD recoveries were within 80 - 120 with the following exceptions: None.

All LCS/LCSD RPDs were less than 20% with the following exceptions: None.

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

SVOA Narration

Were all QA/QC performance criteria specified in the MADEP document CAM achieved? No.

QC Batch 522813 (Samples: CF51974, CF51975, CF51977): -----

The LCS and/or the LCSD recovery is above the upper range for one or more analytes that were not reported in the sample(s), therefore no significant bias is suspected. (2-Nitroaniline)

Instrument:

CHEM07 03/19/20-1

Wes Bryon, Chemist 03/19/20

CF51974 (1X), CF51975 (1X), CF51977 (1X)

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

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

Initial Calibration Evaluation (CHEM07/7_SPLIT_0304):

100% of target compounds met criteria.

The following compounds had %RSDs >20%: None.

The following compounds did not meet recommended response factors: 2-Nitrophenol 0.078 (0.1)

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

Continuing Calibration Verification (CHEM07/0319_06-7_SPLIT_0304) (MCP Compliance):



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MCP Certification Report

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SVOA Narration

Internal standard areas were within 50 to 200% of the initial calibration with the following exceptions: None.

100% of target compounds met criteria.

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

The following compounds did not meet maximum % deviations: None.

The following compounds did not meet recommended response factors: 2-Nitrophenol 0.093 (0.1)

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

QC (Batch Specific):

Batch 522813 (CF51012)

CF51974, CF51975, CF51977

All LCS recoveries were within 40 - 140 with the following exceptions: None.

All LCSD recoveries were within 40 - 140 with the following exceptions: 2-Nitroaniline(147%)

All LCS/LCSD RPDs were less than 20% with the following exceptions: None.

Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid

surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)

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

SVOASIM Narration



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MCP Certification Report

March 24, 2020 SDG I.D.: GCF51973

SVOASIM Narration

Were all QA/QC performance criteria specified in the MADEP document CAM achieved? No.

QC Batch 522813 (Samples: CF51974, CF51975, CF51977): -----

The LCS and/or the LCSD recovery is below the method criteria. All of the other QC is acceptable, therefore no significant bias is suspected. (Hexachlorocyclopentadiene)

The LCS/LCSD RPD exceeds the method criteria for one or more analytes, but these analytes were not reported in the sample(s) so no variability is suspected. (Pyridine)

QC Batch 522853 (Samples: CF51974, CF51975, CF51977): -----

The LCS and/or the LCSD recovery is above the upper range for one or more analytes that were not reported in the sample(s), therefore no significant bias is suspected. (Acenaphthene, Acenaphthylene, Anthracene, Benzo(a)pyrene, Benzo(b)fluoranthene, Benzo(ghi)perylene, Benzo(k)fluoranthene, Chrysene, Dibenz(a,h)anthracene, Fluoranthene, Fluorene, Indeno(1,2,3-cd)pyrene, Phenanthrene)

The LCS and/or the LCSD recovery is above the upper range, therefore a slight high bias is possible. (Pyrene)

The LCS and/or the LCSD recovery is below the method criteria. All of the other QC is acceptable, therefore no significant bias is suspected. (2-Methylnaphthalene)

The LCS/LCSD RPD exceeds the method criteria for one or more analytes, but these analytes were not reported in the sample(s) so no variability is suspected. (Acenaphthene, Acenaphthylene, Anthracene, Benz(a)anthracene, Benzo(b)fluoranthene, Benzo(ghi)perylene, Benzo(k)fluoranthene, Chrysene, Dibenz(a,h)anthracene, Fluoranthene, Fluorene, Indeno(1,2,3-cd)pyrene, Phenanthrene)

The LCS/LCSD RPD exceeds the method criteria for one or more analytes, therefore there may be variability in the reported result. (2-Methylnaphthalene, Naphthalene, Pyrene)

Instrument:

CHEM27 03/18/20-1 Wes Bryon, Chemist 03/18/20

CF51974 (1X), CF51975 (1X), CF51977 (1X)

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

Initial Calibration Evaluation (CHEM27/27_SIM18_0311):

100% of target compounds met criteria.

The following compounds had %RSDs >20%: None.

The following compounds did not meet recommended response factors: None.

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

Continuing Calibration Verification (CHEM27/0318_03-27_SIM18_0311) (MCP Compliance):

Internal standard areas were within 50 to 200% of the initial calibration with the following exceptions: None.

97% of target compounds met criteria.

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

The following compounds did not meet maximum % deviations: None.

The following compounds did not meet recommended response factors: None.



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MCP Certification Report

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SVOASIM Narration

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

QC (Batch Specific):

Batch 522813 (CF51012)

CF51974, CF51975, CF51977

All LCS recoveries were within 40 - 140 with the following exceptions: Hexachlorocyclopentadiene(39%)

All LCSD recoveries were within 40 - 140 with the following exceptions: None.

All LCS/LCSD RPDs were less than 20% with the following exceptions: Pyridine(22.0%)

Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)

Batch 522853 (CF52086)

CF51974, CF51975, CF51977

All LCS recoveries were within 40 - 140 with the following exceptions: Acenaphthene(170%), Acenaphthylene(166%),

Anthracene(158%), Benzo(a)anthracene(175%), Benzo(a)pyrene(163%), Benzo(b)fluoranthene(159%),

Benzo(ghi)perylene(185%), Benzo(k)fluoranthene(149%), Chrysene(165%), Dibenz(a,h)anthracene(171%), Fluoranthene(155%),

Fluorene(165%), Indeno(1,2,3-cd)pyrene(168%), Phenanthrene(151%), Pyrene(163%)

All LCSD recoveries were within 40 - 140 with the following exceptions: 2-Methylnaphthalene(29%)

All LCS/LCSD RPDs were less than 20% with the following exceptions: 2-Methylnaphthalene(77.9%), Acenaphthene(81.0%),

Acenaphthylene(80.2%), Anthracene(80.9%), Benz(a)anthracene(82.3%), Benzo(a)pyrene(81.0%),

 $Benzo(b) fluoranthene (78.9\%), \ Benzo(ghi) perylene (78.2\%), \ Benzo(k) fluoranthene (81.1\%), \ Chrysene (80.9\%), \ Chrysen$

Dibenz(a,h)anthracene(80.3%), Fluoranthene(83.1%), Fluorene(78.5%), Indeno(1,2,3-cd)pyrene(81.2%), Naphthalene(78.5%),

Phenanthrene(82.2%), Pyrene(83.5%)

Additional 8270 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is at least 10%. (Acid surrogates acceptance range for aqueous samples: 10-110%, for soils 30-130%)

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

VOA Narration

Were all QA/QC performance criteria specified in the MADEP document CAM achieved? No.

QC Batch 522985 (Samples: CF51973, CF51974, CF51975, CF51976, CF51977, CF51978, CF51979): -----

The QC recovery for one or more analytes is above the upper range but were not reported in the sample(s), therefore no significant bias is suspected. (trans-1,4-dichloro-2-butene)

QC Batch 523354 (Samples: CF51975): -----

The LCS and/or the LCSD recovery is below the method criteria. All of the other QC is acceptable, therefore no significant bias is suspected. (Trichlorotrifluoroethane)

Instrument:

CHEM02 03/19/20-1

Michael Hahn, Chemist 03/19/20

CF51975 (20X)

Chem02 is a 25ml purge instrument. The laboratory minimum response factor is set at 0.01 instead of 0.05 for the 25ml purge



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MCP Certification Report

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VOA Narration

instruments.

EPA method 8260D Table 4 supports this approach.

Initial Calibration Evaluation (CHEM02/VT-P031720):

98% of target compounds met criteria.

The following compounds had %RSDs >20%: None.

The following compounds did not meet Table 4 recommended minimum response factors: None.

The following compounds did not meet the minimum response factor of 0.05: None.

Continuing Calibration Verification (CHEM02/0319_02-VT-P031720) (MCP Compliance):

Internal standard areas were within 50 to 200% of the initial calibration with the following exceptions: None.

98% of target compounds met criteria.

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

The following compounds did not meet maximum % deviations: None.

The following compounds did not meet Table 4 recommended minimum response factors: None.

The following compounds did not meet the minimum MCP response factor of 0.05: None.

CHEM17 03/17/20-2

Michael Hahn, Chemist 03/17/20

CF51973 (1X), CF51974 (1X), CF51975 (1X), CF51976 (1X), CF51977 (1X), CF51978 (1X), CF51979 (1X)

Chem 17 is a 25ml purge instrument. The laboratory minimum response factor is set at 0.01 instead of 0.05 for the 25ml purge instruments.

EPA method 8260D Table 4 supports this approach.

Initial Calibration Evaluation (CHEM17/VT-S031520):

94% of target compounds met criteria.

The following compounds had %RSDs >20%: 1,2-Dibromo-3-chloropropane 22% (20%), 1,4-Dioxane 28% (20%), Bromomethane 24% (20%), Tetrahydrofuran (THF) 25% (20%)

The following compounds did not meet Table 4 recommended minimum response factors: 1,2-Dibromo-3-chloropropane 0.042 (0.05), Acetone 0.054 (0.1), Bromoform 0.085 (0.1), Methyl ethyl ketone 0.087 (0.1), Tetrahydrofuran (THF) 0.049 (0.05) The following compounds did not meet the minimum response factor of 0.05: 1,2-Dibromo-3-chloropropane 0.042 (0.05),

Tetrahydrofuran (THF) 0.049 (0.05)

Continuing Calibration Verification (CHEM17/0317_24-VT-S031520) (MCP Compliance):

Internal standard areas were within 50 to 200% of the initial calibration with the following exceptions: None.

98% of target compounds met criteria.

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

The following compounds did not meet maximum % deviations: None.

The following compounds did not meet Table 4 recommended minimum response factors: 1,2-Dibromo-3-chloropropane 0.047 (0.05), Acetone 0.049 (0.1), Methyl ethyl ketone 0.081 (0.1), Tetrahydrofuran (THF) 0.042 (0.05)

The following compounds did not meet the minimum MCP response factor of 0.05: 1,2-Dibromo-3-chloropropane 0.042 (0.05), Acetone 0.054 (0.05), Tetrahydrofuran (THF) 0.049 (0.05)

CHEM17 03/18/20-1 Michael Hahn, Chemist 03/18/20

CF51974 (5X), CF51975 (5X), CF51977 (5X)

Chem 17 is a 25ml purge instrument. The laboratory minimum response factor is set at 0.01 instead of 0.05 for the 25ml purge instruments.

EPA method 8260D Table 4 supports this approach.

Initial Calibration Evaluation (CHEM17/VT-S031520):

94% of target compounds met criteria.



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VOA Narration

The following compounds had %RSDs >20%: None.

The following compounds did not meet Table 4 recommended minimum response factors: None.

The following compounds did not meet the minimum response factor of 0.05: None.

Continuing Calibration Verification (CHEM17/0318_02-VT-S031520) (MCP Compliance):

Internal standard areas were within 50 to 200% of the initial calibration with the following exceptions: None.

96% of target compounds met criteria.

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

The following compounds did not meet maximum % deviations: None.

The following compounds did not meet Table 4 recommended minimum response factors: None.

The following compounds did not meet the minimum MCP response factor of 0.05: None.

QC (Batch Specific):

Batch 522985 (CF51399) CHEM17 3/17/2020-2

CF51973(1X), CF51974(1X), CF51975(1X), CF51976(1X), CF51977(1X), CF51978(1X), CF51979(1X)

All LCS recoveries were within 70 - 130 with the following exceptions: trans-1,4-dichloro-2-butene(149%)

All LCSD recoveries were within 70 - 130 with the following exceptions: trans-1,4-dichloro-2-butene(165%)

All LCS/LCSD RPDs were less than 30% with the following exceptions: None.

Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%.

Batch 523171 (CF52205) CHEM17 3/18/2020-1

CF51974(5X), CF51975(5X), CF51977(5X)

All LCS recoveries were within 70 - 130 with the following exceptions: None.

All LCSD recoveries were within 70 - 130 with the following exceptions: None.

All LCS/LCSD RPDs were less than 30% with the following exceptions: None.

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

Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%.

Batch 523354 (CF52176) CHEM02 3/19/2020-1

CF51975(20X)

All LCS recoveries were within 70 - 130 with the following exceptions: Trichlorotrifluoroethane(62%)

All LCSD recoveries were within 70 - 130 with the following exceptions: None.

All LCS/LCSD RPDs were less than 30% with the following exceptions: None.

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

Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%.

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

VPH Narration

Were all QA/QC performance criteria specified in the MADEP document CAM achieved? Yes.

Instrument:

PIDFID 03/18/20-1 Raman Makol, Chemist 03/18/20

CF51974 (1X), CF51975 (1X), CF51977 (1X)



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



MCP Certification Report

March 24, 2020 SDG I.D.: GCF51973

VPH Narration

QC (Batch Specific):

Batch 523130 (CF51974)

CF51974(1X), CF51975(1X), CF51977(1X)

All LCS recoveries were within 70 - 130 with the following exceptions: None.

All LCSD recoveries were within 70 - 130 with the following exceptions: None.

All LCS/LCSD RPDs were less than 20% with the following exceptions: None.

A blank MS/MSD was analyzed with this batch.

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

Environmental Labo Customer: BLCe Address: 3555 Matrix Code: DW-Drinking Water GW-Ground Valve Reaw Water SE-Sediment SL B-Bulk L-Liquid X = SANPLE # MW-10 51975 MW-10 51975 MW-10 51977 MW-10	Coolant: IPK TTCE IN No Terms CHAIN OF CUSTODY RECORD	Data Delivery/Contact Ontion	587 East Middle Tumpike, P.O. Box 370, Manchester, CT 06040	Email: info@phoenixlabs.com Fax (860) 645-0823	Project 1/06 4/2 Broiset D.	PKWK Report to: AMANICAMONING	La CT O Invoice to: Projection BL. Combanies, com	QUOTE# C Bottle Quantities.		ate: 3/17 Analysis Analysis	is anhali	DW=Drinking Warter GW=Ground Warter SW=Surface Warter WW=Waste Water SP=Sediment SL=Sediment SL=Sedime	30 1 10 10 10 10 10 10 10 10 10 10 10 10	Sample Date Time Time Matrix Sampled Sampled Sampled	×	X X X X X X X X X X X X X X X X X X X				> >	×				Sertification	GW Protection	GA Leachability GA Mobility GA CAN 1 CA CAN 2 CA	2.00 2.0 2.00 2.0	Residential DEC
				Environmental Laboratories, Inc.	3	8	ناف		Client Sargele - Information - Identification	With MAG		er GW=Ground Water SW=Surface Wate E=Sediment SL=Sludge S=Soil SD=So (=		Customer Sample Identification	1	MW-2 /	MW-101	MW-102	MW-103	MW-104	Trip Blank		7	Accepted by:	* Kynoller		Comments, Special Requirements or Regulations:		



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

Analysis Report

August 25, 2020

FOR: Attn: Mr. Matt Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Sample Information Custody Information Date <u>Time</u> SURFACE WATER Collected by: 08/21/20 11:30 Matrix: MG Received by: Location Code: **BLCOMPANIES** LB 08/21/20 13:36

Rush Request: 24 Hour Analyzed by: see "By" below

P.O.#: Laboratory Data

SDG ID: GCG59046

Phoenix ID: CG59046

Project ID: 16D3253 Client ID: RW-1

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	08/22/20	CPP	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	08/22/20	CPP	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	08/22/20	CPP	SW6010D
Chromium	< 0.001	0.001	mg/L	1	08/22/20	EK	SW6010D
Copper	< 0.005	0.005	mg/L	1	08/22/20	CPP	SW6010D
Iron	1.80	0.010	mg/L	1	08/22/20	CPP	E200.7
Hardness (CaCO3)	185	0.1	mg/L	1	08/25/20		E200.7
Mercury	< 0.0002	0.0002	mg/L	1	08/24/20	RS	SW7470A
Nickel	0.001	0.001	mg/L	1	08/22/20	CPP	SW6010D
Lead	< 0.002	0.002	mg/L	1	08/22/20	CPP	SW6010D
Antimony	< 0.005	0.005	mg/L	1	08/22/20	CPP	SW6010D
Selenium	< 0.010	0.010	mg/L	1	08/22/20	CPP	SW6010D
Trivalent Chromium	< 0.001	0.001	mg/L	1	08/25/20		Calculation
Zinc	0.004	0.004	mg/L	1	08/22/20	CPP	SW6010D
Chromium, Hexavalent	< 0.01	0.01	mg/L	1	08/24/20 18:19	0	SM3500CRB-11
Ammonia as Nitrogen	0.24	0.05	mg/L	1	08/23/20	ARG	E350.1
рН	7.31	1.00	pH Units	1	08/21/20 23:01	AP/EG	SM4500-H B-11
Salinity	0.5	0.5	ppt	1	08/21/20	AP	SM2520B-10
Mercury Digestion	Completed				08/24/20	VT/VT	SW7470A
Total Metals Digestion	Completed				08/21/20	AG	

Ver 1 Page 1 of 3

Project ID: 16D3253 Phoenix I.D.: CG59046

Client ID: RW-1

RL/

Parameter Result PQL Units Dilution Date/Time By

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

The regulatory hold time for pH is immediately. This pH was performed in the laboratory and may be considered outside of hold-time

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

Phyllis Shiller, Laboratory Director

August 25, 2020

Official Report Release To Follow

Ver 1 Page 2 of 3



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

Analysis Report

August 25, 2020

FOR: Attn: Mr. Matt Gonsalves

BL Companies, Inc. 355 Research Parkway Meriden, CT 06450

Sample Informat	<u>ion</u>	Custody Informat	<u>ion</u>	<u>Date</u>	<u>Time</u>
Matrix:	SURFACE WATER	Collected by:	MG	08/21/20	11:45
Location Code:	BLCOMPANIES	Received by:	LB	08/21/20	13:36

Rush Request: 24 Hour Analyzed by: see "By" below

Laboratory Data

SDG ID: GCG59046

Phoenix ID: CG59047

Project ID: 16D3253 Client ID: MW-2

P.O.#:

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Chromium	< 0.001	0.001	mg/L	1	08/25/20	EK	SW6010D
Trivalent Chromium	< 0.001	0.001	mg/L	1	08/25/20		Calculation
Chlorine Residual	< 0.02	0.02	mg/L	1	08/21/20 18:51	0	SM4500Cl-G-00
Chromium, Hexavalent	< 0.01	0.01	mg/L	1	08/21/20 18:13	0	SM3500CRB-11
Ammonia as Nitrogen	0.27	0.05	mg/L	1	08/23/20	ARG	E350.1
Total Metals Digestion	Completed				08/24/20	AG	

RL/PQL=Reporting/Practical Quantitation Level ND=Not Detected BRL=Below Reporting Level

Comments:

The regulatory hold time for Chlorine is immediately. This Chlorine was performed in the laboratory and may be considered outside of hold-time

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

Phyllis Shiller, Laboratory Director

August 25, 2020

Official Report Release To Follow

Ver 1 Page 3 of 3

Tuesday, August 25, 2020

Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report

GCG59046 - BLCOMPANIES

State: MA

RLAnalysis SampNo Acode Phoenix Analyte Criteria Result RLCriteria Criteria Units

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.

Page 1 of 1

^{***} No Data to Display ***

		2,5						
Cooler: Yes No I I I I I I I I I I I I I I I I I I	This section MUST be completed with Bottle Quantities.		MO SISTED TO STATE OF THE STATE				Data Format Excel GIS/Key GIS/Key Other Data Package Tier II Checklist Full Data Package* Dhoenix Std Report Other	* SURCHARGE APPLIES
Coolen: PK Yes Coolen: PK N ICE Temp 0, 4° C Pg o Data Delivery/Contact Options: Fax: Phone: Email: Project DO:			OT WO		m		MA MCP Certification MCP Certification CW-1 MWRA esMART GW-2 GW-2 S-1 GW-3 S-2 GW-1 S-2 GW-2 S-2 GW-3 S-3 GW-1 S-3 GW-2 S-3 GW-3 S-3 GW-1 S-3 GW-2 S-3 GW-3 MA MCP M	
JRD ster, CT 06040 645-0823 26	& Bloom				X		MA MCP Certification	-
East Middle Turnpike, P.O. Box 370, Manchester, CT 06040 Email: info@phoenixlabs.com Fax (860) 645-0823 Client Services (860) 645-8726 Droingt: 17.0.23.673	mesossky			H	×		oility	Objectives
CHAIN OF CUSTC Last Middle Tumpike, P.O. Boy Email: info@phoenixlabs.com Client Services (8	Project: Report to: Invoice to: QUOTE#	Analysis Request		^ ×			17 3 3 C	SURCHARGE APPLIES
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PHOEINITY SET OF THE STATE OF T	Meridea CT	Sampler's Signature MAR G. Date: 8/2 Matrix Code: DW=Drinking Water GW=Ground Water SW=Surface Water WW=Waste Water SR=Sediment SL=Sludge S=Soil SD=Solid W=Wipe OIL=Oil	Customer Sample Identification	RW-1	MW-2		Chromium L. Coper, Lead, Merch C. Nickel, Selevium, Silver, 2:00	you winds
PHO Environmen	Address: Address: A	Sampler's Signature Matrix Code: DW=Drinking Water RW=Raw Water SE B=Bulk L=Liquid X=	PHOENIX USE ONLY SAMPLE #	59046	C4047		Comments, Special F Chromium Chromium Vicke Se	1.11010 761 7

GCG 59046

Shannon Wilhelm

Subject:

FW: 16D3253

From: Gonsalves, Matthew [mailto:mgonsalves@Blcompanies.com]

Sent: Monday, August 24, 2020 12:31 PM

To: Christine Paradise

Subject: Re: 16D3253

Yes please proceed with running them as soon as possible.

Get Outlook for Android

From: Christine Paradise <christine@phoenixlabs.com>

Sent: Monday, August 24, 2020 12:11:42 PM

To: Gonsalves, Matthew <mgonsalves@Blcompanies.com>

Subject: 16D3253

Good Afternoon,

We received these samples on Friday, however the HexCr and TriCr was not logged into the first sample due to our error. Would you like to proceed with testing past holding time or is this something you would need to resample? Please let me know.

I'm sorry for any inconvenience.

Christine Paradise

Phoenix Environmental Laboratories

587 East Middle Turnpike

Manchester, CT 06040

Phone: 860-645-1102 Fax: 860-645-0823

ATTACHMENT G
TREATMENT
SYSTEM
SPECIFICATIONS
AND DATA SHEETS



CGS

CATION EXCHANGE RESIN SOFTENING GRADE Na FORM

RESINTECH CGS is a sodium form standard crosslinked gel strong acid cation resin. *CGS* is optimized for residential applications that require good regeneration efficiency and high capacity. *RESINTECH CGS* is intended for use in all residential and commercial softening applications that do not have significant amounts of chlorine in the feedwater. *CGS* is supplied in the sodium form.



FEATURES & BENEFITS

RESIDENTIAL SOFTENING APPLICATIONS

Resin parameters are optimized for residential softeners

LOW COLOR THROW

SUPERIOR PHYSICAL STABILITY

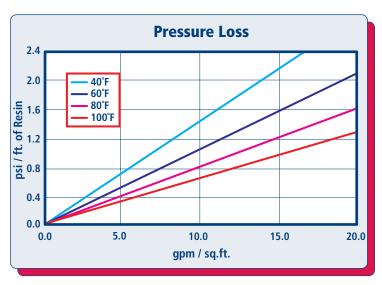
93% plus sphericity and high crush strengths together with carefully controlled particle distribution provides long life and low pressure drop

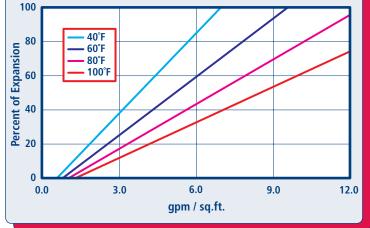
COMPLIES WITH US FDA REGULATIONS

Conforms to paragraph 21CFR173.25 of the Food Additives Regulations of the US FDA

Prior to first use for potable water, resin should be backwashed for a minimum of 20 minutes, followed by 10 bed volumes of downflow rinse.

HYDRAULIC PROPERTIES





Backwash Expansion

PRESSURE LOSS

The graph above shows the expected pressure loss of *ResinTech CGS* per foot of bed depth as a function of flow rate at various temperatures.

BACKWASH

The graph above shows the expansion characteristics of *ResinTech CGS* as a function of flow rate at various temperatures.

RESINTECH CGS

PHYSICAL PROPERTIES

Polymer Structure Styrene/DVB

Polymer Type Gel

Functional Group Sulfonic Acid Physical Form Spherical beads

Ionic Form as shipped Sodium

Total Capacity

Sodium form >1.8 meg/mL

Water Retention

Sodium form 40 to 52 percent

Approximate Shipping Weight

Sodium form 50 lbs./cu.ft.

Screen Size Distribution (U.S. mesh) 16 to 50

Maximum Fines Content (<50 mesh) 1 percent

Minimum Sphericity 90 percent

Uniformity Coefficient 1.6 approx.

Resin Color Amber

Note: Physical properties can be certified on a per lot basis, available upon request

SUGGESTED OPERATING CONDITIONS

Maximum continuous temperature

Sodium form 250°F Minimum bed depth 24 inches

Backwash expansion 25 to 50 percent

Maximum pressure loss 25 psi
Operating pH range 0 to 14 SU

Regenerant Concentration

Salt cycle 10 to 15 percent NaCl Regenerant level 4 to 15 lbs./cu.ft. Regenerant flow rate. 0.5 to 1.5 gpm/cu.ft.

Regenerant contact time >20 minutes

Displacement flow rate

Displacement volume

10 to 15 gallons/cu.ft.

Rinse flow rate

Same as service flow

Rinse volume

35 to 60 gallons/cu.ft.

Service flow rate

1 to 10 gpm/cu.ft.

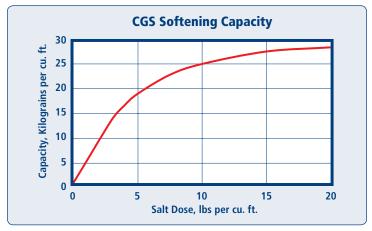
Note: These guidelines describe average low risk operating conditions. They are not intended to be absolute minimums or maximums.

For operation outside these guidelines, contact ResinTech Technical Support

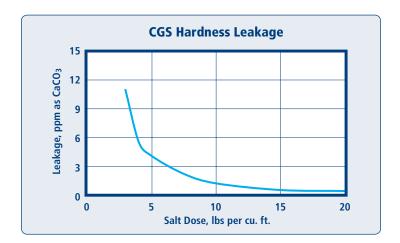
APPLICATIONS

SOFTENING

RESINTECH CGS is a standard crosslinked cation resin optimized for residential and commercial applications. This type of resin is easier to regenerate than the higher crosslinked resins. CGS has marginal resistance to chlorine and other oxidants and is not ideal for high temperature and other high stress applications.



Capacity and leakage data are based on the following: 2:1 Ca:Mg ratio, 500 ppm TDS as CaCO3, 0.2% hardness in the salt and 10% brine concentration applied co-currently through the resin over 30 minutes. No engineering downgrade has been applied.





East Coast - West Berlin, NJ p:856.768.9600 • Midwest - Chicago, IL p:708.777.1167 • West Coast - Los Angeles, CA p:323.262.1600



89 Crawford Street

Leominster, Massachusetts 01453

Tel: 774.450.7177 Fax: 888.835.0617 www.lrt-llc.net

FILTRATION MEDIA: 8x30 RE-ACTIVATED CARBON 4x10 RE-ACTIVATED CARBON

GENERAL DESCRIPTION

Select Re-Activated carbon from domestic sources is quality screened during our purchasing process for activity, density and fines. The use of re-activated carbon is recommended as a lower cost alternative for most sites where drinking water quality is not necessary. In many cases our re-activated carbon meets and exceeds imported virgin carbon. In addition all carbon either sold by itself or installed in our filtration units traced by lot number to the installation or sale.

8x30 (Liquid Phase) Standard Specifications:	Standard	Value
lodine Number	ASTM D-4607	800 Minimum
Moisture Content	ASTM D-2867	5% Maximum (as packed)
Particle Size	ASTM D-2862	8x30 US Mesh
Ash		10% Maximum
Total Surface Area (N2BET)		1050 Minimum
Pore Volume (cc/g)		0.75

4*10 (Vapor Phase) Standard Specifications:	Standard	Value
Carbon Tetrachloride Activity Level	ASTM D-3467	40 Minimum
Moisture Content	ASTM D-2867	5% Maximum (as packed)
Particle Size	ASTM D-2862	4x10 US Mesh
Ash		10% Maximum
Total Surface Area (N2BET)		1050 Minimum
Pore Volume (cc/g)		0.75



The Pulsatron Series HV designed for high viscosity applications for precise and accurate metering control. The Series HV offers manual control over stroke length and stroke rate as standard with the option to choose between 4-20mA and external pace inputs for automatic control.

Five distinct models are available, having pressure capabilities to 150 PSIG (10 BAR) @ 12 GPD (1.9 lph), and flow capacities to 240 GPD (37.9 lph) @ 80 PSIG (5.6 BAR), with a turndown ratio of 100:1. Metering performance is reproducible to within ± 2% of maximum capacity.

Features

- Automatic Control, available with 4-20mADC direct or external pacing, with stop function.
- Manual Control by on-line adjustable stroke rate and stroke length.
- Auto-Off-Manual switch.
- · Highly Reliable timing circuit.
- · Circuit Protection against voltage and current upsets.
- Panel Mounted Fuse.
- Solenoid Protection by thermal overload with autoreset
- Water Resistant, for outdoor and indoor applications.
- Indicator Lights, panel mounted.
- Guided Ball Check Valve Systems, to reduce back flow and enhance outstanding priming characteristics.
- Viscosities to 20,000 CPS.

Controls



Manual Stroke Rate

• Turn-Down Ratio 10:1

Manual Stroke Length

Turn-Down Ratio 10:1

4-20mA or 20-4mA Input

Automatic Control

Operating Benefits

- Reliable metering performance.
- Rated "hot" for continuous duty.
- High viscosity capability.
- Leak-free, sealless, liquid end.



Aftermarket

- KOPkits
- Gauges
- Dampeners
- Pressure Relief Valves
- Tanks
- Pre-Engineered Systems
- Process Controllers (PULSAblue, MicroVision)











Series HV

Specifications and Model Selection

MODEL		LVB3	LVF4	LVG4	LVG5	LVH7
Capacity	GPH	0.50	1.00	2.00	4.00	10.00
nominal	GPD	12	24	48	96	240
(max.)	LPH	1.9	3.8	7.6	15.1	37.9
Pressure	PSIG	150	150	110	110	80
(max.)	BAR	10	10	7	7	5.6
Connections:	Tubing	STATE OF STREET		38" I.D. X .5		SECURE OF THE PROPERTY.



Engineering Data

Pump Head Materials Available: GFPPL

PVC PVDF 316 SS

Diaphragm: PTFE-faced CSPE-backed

Check Valves Materials Available:

Seats/O-Rings: PTFE

CSPE Viton

Balls: Ceramic

PTFE 316 SS Alloy C GFPPL

Fittings Materials Available: GFPI
PVC

PVC PVDF

Bleed Valve: Same as fitting and check valve

selected, except 316SS

Injection Valve & Foot Valve Assy: Same as fitting and check valve

selected

Tubing: Clear PVC White PE

Important: Material Code - GFPPL=Glass-filled Polypropylene, PVC=Polyvinyl Chloride, PE=Polyethylene, PVDF=Polyvinylidene Fluoride, CSPE=Generic formulation of Hypalon, a registered trademark of E.I. DuPont Company. Viton is a registered trademark of E.I. DuPont Company. PVC wetted end recommended for sodium hypochlorite.

Engineering Data

Reproducibility: +/- 2% at maximum capacity

Viscosity Max CPS: 20,000 CPS

Stroke Frequency Max SPM:125Stroke Frequency Tum-Down Ratio:10:1Stroke Length Tum-Down Ratio:10:1

Power Input: 115 VAC/50-60 HZ/1 ph 230 VAC/50-60 HZ/1 ph

Average Current Draw:

@ 115 VAC; Amps: 1.0 Amps

@ 230 VAC; Amps: 0.5 Amps @ 230 VAC

Peak Input Power: 300 Watts Average Input Power @ Max SPM: 130 Watts

Custom Engineered Designs – Pre-Engineered Systems



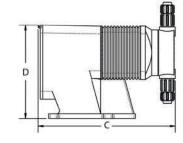
Pre-Engineered Systems

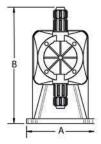
Pulsafeeder's Pre-Engineered Systems are designed to provide complete chemical feed solutions for all electronic metering applications. From stand alone simplex pH control applications to full-featured, redundant sodium hypochlorite disinfection metering, these rugged fabricated assemblies offer tum-key simplicity and industrial-grade durability. The UV-stabilized, high-grade HDPE frame offers maximum chemical compatibility and structural rigidity. Each system is factory assembled and hydrostatically tested prior to shipment.

Dimensions

Series HV Dimensions (inches)											
Model No.	Α	В	С	D	Shipping Weight						
LVB3	5.4	9.3	9.5	7.5	13						
LVF4	5.4	10.8	10.8	7.5	18						
LVG4	5.4	9.5	10.6	7.5	18						
LVG5	5.4	10.8	10.8	7.5	18						
LVH7	6.1	11.5	11	8.2	25						

NOTE: Inches X 2.54 = cm

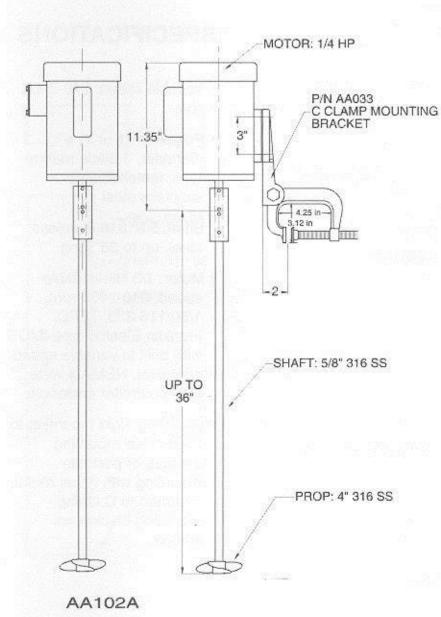








MIXER MODEL NO. AA102A



SPECIFICATIONS

- Speed: 1,725 rpm
- Propeller: (1 or 2)
 4" diameter, 3 blade marine type, material: 316 stainless steel
- Shaft: 5/8" 316 stainless steel, up to 36" long
- Motor: 1/4 HP, 1,725 rpm, 1/60/115-230, capacitor start, or 3/60/230-460, TEFC
- Mounting: rigid mounting to fixed mixer mounting bracket, or portable mounting with mixer motor mounted to C clamp mounting bracket no. AA033.



Revision date 2019-15-4

SAFETY DATA SHEET

Revision number 1

SECTION 1) CHEMICAL PRODUCT AND SUPPLIER'S IDENTIFICATION

Product Name: Redux E50

Product Use: Water and Wastewater Treatment Coagulant/Flocculant

Revision Date: Apr 15, 2019
Supersedes Date: Mar 5, 2015

Manufacturer's Name: Azure Water Services

Address: 280 Callegari Dr. West Haven CT, 06516

Emergency Phone: Chemtrec, (1) 800-424-9300, in US and Canada only

SECTION 2) HAZARDS IDENTIFICATION

Classification

Corrosive to metals - Category 1

Eye Irritation - Category 2

Skin Irritation - Category 2

Pictograms



Signal Word

Warning

Hazardous Statements - Health

Causes serious eye irritation

Causes skin irritation

Hazardous Statements - Physical

May be corrosive to metals

Precautionary Statements - General

If medical advice is needed, have product container or label at hand.

Keep out of reach of children.

Read label before use.

Precautionary Statements - Prevention

Keep only in original packaging.

Wash thoroughly after handling.

Wear protective gloves/protective clothing/eye protection/face protection.



Precautionary Statements - Response

Absorb spillage to prevent material damage.

IF IN EYES: Rinse cautiously with water for several minutes. Remove contact lenses, if present and easy to do. Continue rinsing.

If eye irritation persists: Get medical advice/attention.

IF ON SKIN: Wash with plenty of water.

Specific treatment (see first-aid on this SDS).

If skin irritation occurs: Get medical advice/attention.

Take off contaminated clothing. And wash it before reuse.

Precautionary Statements - Storage

Store in a corrosive resistant container with a resistant inner liner.

Precautionary Statements - Disposal

No precautionary statement available.

Hazards Not Otherwise Classified (HNOC)

None.

SECTION 3) COMPOSITION / INFORMATION ON INGREDIENTS

CAS Chemical Name % By Weight
PROPRIETARY Trade Secret Ingredient 45 - 55%

Specific chemical identity and/or exact percentage (concentration) of the composition has been withheld to protect confidentiality.

SECTION 4) FIRST-AID MEASURES

Inhalation

Remove source of exposure or move person to fresh air and keep comfortable for breathing. Immediately call a POISON CENTER/doctor/. If breathing has stopped, trained personnel should begin rescue breathing or, if the heart has stopped, immediately start cardiopulmonary resuscitation (CPR) or automated external defibrillation (AED).

Eye Contact

Remove source of exposure or move person to fresh air. Rinse eyes cautiously with lukewarm, gently flowing water for several minutes, while holding the eyelids open. Remove contact lenses, if present and easy to do. Continue rinsing for a flushing duration of 30 minutes. Take care not to rinse contaminated water into the unaffected eye or onto the face. Immediately call a POISON CENTER/doctor.

Skin Contact

Take off immediately all contaminated clothing, shoes and leather goods (e.g. watchbands, belts). Rinse skin with lukewarm, gently flowing water/shower for a duration of 30 minutes or until medical aid is available. Immediately call a POISON CENTER/doctor. Wash contaminated clothing before re-use or discard.

Ingestion

Rinse mouth with water. Do NOT induce vomiting. Give 1 to 2 cups of milk or water to drink. Never give anything by mouth to an unconscious person. If vomiting occurs naturally, lie on your side, in the recovery position. Immediately call a POISON CENTER/doctor.

Most Important Symptoms and Effects, Both acute and Delayed

No data available.

Indication of Any Immediate Medical Attention and Special Treatment Needed

No data available.

SECTION 5) FIRE-FIGHTING MEASURES

Suitable Extinguishing Media

Dry chemical, foam, carbon dioxide. Sand or earth may be used for small fires only.

Use extinguishing agent suitable for type of surrounding fire.

Redux E50 Page 2 of 6

Unsuitable Extinguishing Media

Do not use direct water stream since this may cause fire to spread.

Specific Hazards in Case of Fire

In case of fire, hazardous decomposition products may include sulphur oxides.

Fire-Fighting Procedures

Isolate immediate hazard area and keep unauthorized personnel out. Stop spill/release if it can be done safely. Move undamaged containers from immediate hazard area if it can be done safely. Water spray may be useful in minimizing or dispersing vapors and to protect personnel. Water may be ineffective but can be used to cool containers exposed to heat or flame. Caution should be exercised when using water or foam as frothing may occur, especially if sprayed into containers of hot, burning liquid. Dispose of fire debris and contaminated extinguishing water in accordance with official regulations.

Special Protective Actions

Wear protective pressure self-contained breathing apparatus (SCBA) and full turnout gear.

SECTION 6) ACCIDENTAL RELEASE MEASURES

Emergency Procedure

Isolate hazard area and keep unnecessary people away. Remove all possible sources of ignition in the surrounding area. Notify authorities if any exposure to the general public or the environment occurs or is likely to occur.

Absorb spill with absorbent material or vacuum spill into polyethylene lined steel or plastic drums.

Do not touch or walk through spilled material.

If spilled material is cleaned up using a regulated solvent, the resulting waste mixture may be regulated.

Recommended Equipment

Positive pressure, full-facepiece self-contained breathing apparatus (SCBA), or positive pressure supplied air respirator with escape SCBA (NIOSH approved).

Personal Precautions

Avoid breathing vapor or mist. Avoid contact with skin, eye or clothing. Ensure adequate ventilation. Do not touch damaged containers or spilled materials unless wearing appropriate protective clothing.

Environmental Precautions

Stop spill/release if it can be done safely. Prevent spilled material from entering sewers, storm drains, other unauthorized drainage systems and natural waterways by using sand, earth, or other appropriate barriers.

Methods and Materials for Containment and Cleaning Up

Contain and collect spillage with non-combustible, absorbent material e.g. sand, earth, vermiculite or diatomaceous earth and place in container for disposal according to local regulations. Contaminated absorbent material may pose the same hazard as the spilled product.

SECTION 7) HANDLING AND STORAGE

General

Wash hands after use.

Do not get in eyes, on skin or on clothing.

Do not breathe vapors or mists.

Use good personal hygiene practices.

Eating, drinking and smoking in work areas is prohibited.

Remove contaminated clothing and protective equipment before entering eating areas.

Eyewash stations and showers should be available in areas where this material is used and stored.

Ventilation Requirements

Use only with adequate ventilation to control air contaminants to their exposure limits. The use of local ventilation is recommended to control emissions near the source.

Storage Room Requirements

Keep container(s) tightly closed and properly labeled. Store in cool, dry, well-ventilated areas away from heat, direct sunlight and strong oxidizers. Store in approved containers and protect against physical damage. Keep containers securely sealed when not in use. Indoor storage should meet OSHA standards and appropriate fire codes. Containers that have been opened must be carefully resealed to prevent leakage. Empty containers retain residue and may be dangerous.

Use non-sparking ventilation systems, approved explosion-proof equipment and intrinsically safe electrical systems in areas where this product is used and stored.

Redux E50 Page 3 of 6

SECTION 8) EXPOSURE CONTROLS, PERSONAL PROTECTION

Eye Protection

Wear eye protection with side shields or goggles. Wear indirect-vent, impact and splash resistant goggles when working with liquids. If additional protection is needed for entire face, use in combination with a face shield.

Skin Protection

Use of gloves approved to relevant standards made from the following materials may provide suitable chemical protection: PVC, neoprene or nitrile rubber gloves. Suitability and durability of a glove is dependent on usage, e.g. frequency and duration of contact, chemical resistance of glove material, glove thickness, dexterity. Always seek advice from glove suppliers. Contaminated gloves should be replaced. Use of an apron and over-boots of chemically impervious materials such as neoprene or nitrile rubber is recommended to avoid skin sensitization. The type of protective equipment must be selected according to the concentration and amount of the dangerous substance at the specific workplace. Launder soiled clothes or properly disposed of contaminated material, which cannot be decontaminated.

Respiratory Protection

If engineering controls do not maintain airborne concentrations to a level which is adequate to protect worker, a respiratory protection program that meets or is equivalent to OSHA 29 CFR 1910.134 and ANSI Z88.2 should be followed. Check with respiratory protective equipment suppliers.

Appropriate Engineering Controls

Provide exhaust ventilation or other engineering controls to keep the airborne concentrations of vapors below their respective threshold limit value.

SECTION 9) PHYSICAL AND CHEMICAL PROPERTIES

Physical and Chemical Properties

Density	11.10 lb/gal
Specific Gravity	1.33 - 1.35
Appearance	Colorless to yellow liquid
рН	3 - 4
Odor Threshold	N/A
Odor Description	N/A
Water Solubility	complete
Viscosity	< 100cps @20C
Vapor Pressure	Similar to water
Vapor Density	N/A
Freezing Point	<19 °F
Boiling Point	>212 °F
Evaporation Rate	N/A
Flammability	Will not burn

SECTION 10) STABILITY AND REACTIVITY

Stability

Stable under normal storage and handling conditions.

Conditions To Avoid

Avoid heat, sparks, flame, high temperature and contact with incompatible materials.

Hazardous Reactions/Polymerization

Hazardous polymerization will not occur.

Incompatible Materials

Strong bases, acids, oxidizing and reducing agents.

Hazardous Decomposition Products

May produce carbon monoxide, carbon dioxide.

Redux E50 Page 4 of 6

SECTION 11) TOXICOLOGICAL INFORMATION

Likely Routes of Exposure

Inhalation LC50 : Not Available Oral LD50 : Not Available Dermal LD50 : Not Available

Acute Toxicity

Component	weight-%	Oral LD50	Dermal LD50	Inhalation LC50
Trade Secret Ingredient	45 - 55%	= 9187 mg/kg (Rat)	> 2000 mg/k (Rat)	

Aspiration Hazard

No Data Available

Respiratory/Skin Sensitization

No Data Available

Serious Eye Damage/Irritation

Causes serious eye irritation

Skin Corrosion/Irritation

Causes skin irritation

Specific Target Organ Toxicity - Repeated Exposure

No Data Available

Specific Target Organ Toxicity - Single Exposure

No Data Available

SECTION 12) ECOLOGICAL INFORMATION

Ecotoxicity

Acute aquatic toxicity - Product Information

Fish LC 50 (96 hour, static) 776.4 mg/L Pimephales promelas (Fathead Minnow) 1

EC 50 (96 hour, static) 265.5 mg/L Pimephales promelas (Fathead Minnow) 1

Crustacea LC 50 (48 hour, static) 803.8 mg/L Ceriodaphnia dubia (Water Flea) 1

EC 50 (48 hour, static) 33.2 mg/L Ceriodaphnia dubia (Water Flea)

Algae/aquatic plants No information available

Acute aquatic toxicity - Component Information

Component	weight-%	Algae/aquatic plants	Fish	Toxicity to daphnia and other aquatic invertebrates
Trade Secret Ingredient	45 - 55%	1	LC50 (96 h static) 100 - 500 mg/L (Brachydanio rerio)	

Mobility in Soil

No data available.

Bio-accumulative Potential

No data available.

Persistence and Degradability

No data available.

Other Adverse Effect

No data available.

Redux E50 Page 5 of 6

SECTION 13) DISPOSAL CONSIDERATIONS

Waste Disposal

Under RCRA it is the responsibility of the user of the product to determine at the time of disposal whether the product meets RCRA criteria for hazardous waste. Waste management should be in full compliance with federal, state and local laws. Empty Containers retain product residue which may exhibit hazards of material, therefore do not pressurize, cut, glaze, weld or use for any other purposes. Return drums to reclamation centers for proper cleaning and reuse.

SECTION 14) TRANSPORT INFORMATION

U.S. DOT Information

NOT REGULATED FOR TRANSPORTATION

This product is excepted from DOT regulations under 49 CFR 173.154(d) when shipped by road or railway. The product exception is referenced in 49 CFR 172.101 Table. Packaging material must not be aluminum, steel or be degraded by this product

SECTION 15) REGULATORY INFORMATION

CAS	Chemical Name	% By Weight	Regulation List
No applicable CAS	No applicable chemical	-	-

SECTION 16) OTHER INFORMATION

Glossary

ACGIH- American Conference of Governmental Industrial Hygienists; ANSI- American National Standards Institute; Canadian TDGCanadian Transportation of Dangerous Goods; CAS- Chemical Abstract Service; Chemtrec- Chemical Transportation Emergency Center(US); CHIP- Chemical Hazard Information and Packaging; DSL- Domestic Substances List; EC- Equivalent Concentration; EH40 (UK)- HSE Guidance Note EH40 Occupational Exposure Limits; EPCRA- Emergency Planning and Community Right-To-Know Act; ESL Effects screening levels; HMIS- Hazardous Material Information Service; LC- Lethal Concentration; LD- Lethal Dose; NFPA- National Fire Protection Association; OEL- Occupational Exposure Limits; OSHA- Occupational Safety and Health Administration, US Department of Labor; PEL- Permissible Exposure Limit; SARA (Title III)- Superfund Amendments and Reauthorization Act; SARA 313- Superfund Amendments and Reauthorization Act, Section 313; SCBA- Self Contained Breathing Apparatus; STEL-Short Term Exposure Limit; TCEQ Texas Commission on Environmental Quality; TLV- Threshold Limit Value; TSCA- Toxic Substances Control Act Public Law 94-469; TWA Time Weighted Value; US DOT- US Department of Transportation; WHMIS- Workplace Hazardous Materials Information System.

Additional Information

Any concentration shown as a range is to protect confidentiality or is due to batch variation.

Version 1.0:

Revision Date: Apr 15,2019

First Edition.

DISCLAIMER

To the best of our knowledge, the information contained herein is accurate. However, neither the above named supplier nor any of its subsidiaries assumes any liability whatsoever for the accuracy or completeness of the information contained herein. Final determination of suitability of any material is the sole responsibility of the user. All materials may present unknown hazards and should be used with caution. Although certain hazards are described herein, we cannot guarantee that these are the only hazards that exist. The above information pertains to this product as currently formulated, and is based on the information available at this time. Addition of reducers or other additives to this product may substantially alter the composition and hazards of the product. Since conditions of use are outside our control, we make no warranties, express or implied, and assume no liability in connection with any use of this information.

Redux E50 Page 6 of 6



Revision date 2019-15-4

SAFETY DATA SHEET

Revision number 1

SECTION 1) CHEMICAL PRODUCT AND SUPPLIER'S IDENTIFICATION

Product ID: Redux-823

Product Name: Processing aid for industrial applications

Revision Date: Apr 15, 2019 Supersedes Date: Jan 25, 2018

Manufacturer's Name: Azure Water Services

Address: 280 Callegari Drive West Haven, CT, US, 06516

Emergency Phone: Chemtrec 800-424-9300, in US and Canada only

SECTION 2) HAZARDS IDENTIFICATION

Classification of the substance or mixture

Not a hazardous substance or mixture according to United States Occupational Safety and Health Administration (OSHA) Hazard Communication Standard (29 CFR 1910.1200).

Hazards Not Otherwise Classified (HNOC)

None.

SECTION 3) COMPOSITION / INFORMATION ON INGREDIENTS

None of the chemicals in this product are hazardous according to the GHS.

SECTION 4) FIRST-AID MEASURES

Inhalation

Remove source of exposure or move person to fresh air and keep comfortable for breathing. Immediately call a POISON CENTER/doctor/. If breathing has stopped, trained personnel should begin rescue breathing or, if the heart has stopped, immediately start cardiopulmonary resuscitation (CPR) or automated external defibrillation (AED).

Eye Contact

Remove source of exposure or move person to fresh air. Rinse eyes cautiously with lukewarm, gently flowing water for several minutes, while holding the eyelids open. Remove contact lenses, if present and easy to do. Continue rinsing for a flushing duration of 30 minutes. Take care not to rinse contaminated water into the unaffected eye or onto the face. Immediately call a POISON CENTER/doctor.

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Ingestion

Rinse mouth with water. Do NOT induce vomiting. Give 1 to 2 cups of milk or water to drink. Never give anything by mouth to an unconscious person. If vomiting occurs naturally, lie on your side, in the recovery position. Immediately call a POISON CENTER/doctor.

Most Important Symptoms and Effects, Both acute and Delayed

No data available.

Indication of Any Immediate Medical Attention and Special Treatment Needed

No data available.



SECTION 5) FIRE-FIGHTING MEASURES

Suitable Extinguishing Media

Dry chemical, foam, carbon dioxide. Sand or earth may be used for small fires only.

Use extinguishing agent suitable for type of surrounding fire.

Unsuitable Extinguishing Media

Do not use direct water stream since this may cause fire to spread.

Specific Hazards in Case of Fire

In case of fire, hazardous decomposition products may include sulphur oxides.

Fire-Fighting Procedures

Isolate immediate hazard area and keep unauthorized personnel out. Stop spill/release if it can be done safely. Move undamaged containers from immediate hazard area if it can be done safely. Water spray may be useful in minimizing or dispersing vapors and to protect personnel. Water may be ineffective but can be used to cool containers exposed to heat or flame. Caution should be exercised when using water or foam as frothing may occur, especially if sprayed into containers of hot, burning liquid. Dispose of fire debris and contaminated extinguishing water in accordance with official regulations.

Special Protective Actions

Wear protective pressure self-contained breathing apparatus (SCBA) and full turnout gear.

SECTION 6) ACCIDENTAL RELEASE MEASURES

Emergency Procedure

Isolate hazard area and keep unnecessary people away. Remove all possible sources of ignition in the surrounding area. Notify authorities if any exposure to the general public or the environment occurs or is likely to occur.

Absorb spill with absorbent material or vacuum spill into polyethylene lined steel or plastic drums.

Do not touch or walk through spilled material.

If spilled material is cleaned up using a regulated solvent, the resulting waste mixture may be regulated.

Recommended Equipment

Positive pressure, full-facepiece self-contained breathing apparatus (SCBA), or positive pressure supplied air respirator with escape SCBA (NIOSH approved).

Personal Precautions

Avoid breathing vapor or mist. Avoid contact with skin, eye or clothing. Ensure adequate ventilation. Do not touch damaged containers or spilled materials unless wearing appropriate protective clothing.

Environmental Precautions

Stop spill/release if it can be done safely. Prevent spilled material from entering sewers, storm drains, other unauthorized drainage systems and natural waterways by using sand, earth, or other appropriate barriers.

Methods and Materials for Containment and Cleaning Up

Contain and collect spillage with non-combustible, absorbent material e.g. sand, earth, vermiculite or diatomaceous earth and place in container for disposal according to local regulations. Contaminated absorbent material may pose the same hazard as the spilled product.

SECTION 7) HANDLING AND STORAGE

General

Wash hands after use.

Do not get in eyes, on skin or on clothing.

Do not breathe vapors or mists.

Use good personal hygiene practices.

Eating, drinking and smoking in work areas is prohibited.

Remove contaminated clothing and protective equipment before entering eating areas.

Eyewash stations and showers should be available in areas where this material is used and stored.

Ventilation Requirements

Use only with adequate ventilation to control air contaminants to their exposure limits. The use of local ventilation is recommended to control emissions near the source.

Storage Room Requirements

Keep container(s) tightly closed and properly labeled. Store in cool, dry, well-ventilated areas away from heat, direct sunlight and strong oxidizers. Store in approved containers and protect against physical damage. Keep containers securely sealed when not in use. Indoor storage should meet OSHA standards and appropriate fire codes. Containers that have been opened must be carefully resealed to prevent leakage. Empty containers retain residue and may be dangerous.

Use non-sparking ventilation systems, approved explosion-proof equipment and intrinsically safe electrical systems in areas where this product is used and stored.

SECTION 8) EXPOSURE CONTROLS, PERSONAL PROTECTION

Eye Protection

Wear eye protection with side shields or goggles. Wear indirect-vent, impact and splash resistant goggles when working with liquids. If additional protection is needed for entire face, use in combination with a face shield.

Skin Protection

Use of gloves approved to relevant standards made from the following materials may provide suitable chemical protection: PVC, neoprene or nitrile rubber gloves. Suitability and durability of a glove is dependent on usage, e.g. frequency and duration of contact, chemical resistance of glove material, glove thickness, dexterity. Always seek advice from glove suppliers. Contaminated gloves should be replaced. Use of an apron and over-boots of chemically impervious materials such as neoprene or nitrile rubber is recommended to avoid skin sensitization. The type of protective equipment must be selected according to the concentration and amount of the dangerous substance at the specific workplace. Launder soiled clothes or properly disposed of contaminated material, which cannot be decontaminated.

Respiratory Protection

If engineering controls do not maintain airborne concentrations to a level which is adequate to protect worker, a respiratory protection program that meets or is equivalent to OSHA 29 CFR 1910.134 and ANSI Z88.2 should be followed. Check with respiratory protective equipment suppliers.

Appropriate Engineering Controls

Provide exhaust ventilation or other engineering controls to keep the airborne concentrations of vapors below their respective threshold limit value.

SECTION 9) PHYSICAL AND CHEMICAL PROPERTIES

Physical and Chemical Properties Density 6.26 lb/gal Specific Gravity 0.6 - 0.9Appearance granular, white solid 5 - 9 @ 5 g/L Odor Threshold N/A Odor Description N/A Water Solubility Complete N/A Viscosity Vapor Pressure Similar to water Vapor Density N/A Freezing Point <32 °F **Boiling Point** >212 °F **Evaporation Rate** N/A Flammability Will not burn

Redux-823 Page 3 of 5

SECTION 10) STABILITY AND REACTIVITY

Stability

Stable under normal storage and handling conditions.

Conditions To Avoid

Avoid heat, sparks, flame, high temperature and contact with incompatible materials.

Hazardous Reactions/Polymerization

Hazardous polymerization will not occur.

Incompatible Materials

Strong bases, acids, oxidizing and reducing agents.

Hazardous Decomposition Products

May produce carbon monoxide, carbon dioxide.

SECTION 11) TOXICOLOGICAL INFORMATION

Likely Routes of Exposure

No Data Available

Acute Toxicity

Inhalation, Testing: Not expected to be toxic by inhalation.

Ingestion, Testing: LD50, Rat > 5,00 mg/kg Dermal, Testing: LD50, Rat > 5,000 mg/kg

Respiratory/Skin Sensitization

No Data Available

Serious Eye Damage/Irritation

No Data Available

Skin Corrosion/Irritation

No Data Available

Specific Target Organ Toxicity - Repeated Exposure

No Data Available

Specific Target Organ Toxicity - Single Exposure

No Data Available

SECTION 12) ECOLOGICAL INFORMATION

Acute Ecotoxicity

Danio Rerio: 96 hr LC50 >100 mg/l (OECD 203)

Fathead Minnow (pimephales promelas): 96hr LC50 >100 mg/l (OECD 203)

Daphnia Magna: 48hr EC50 >100 mg/l (OECD 202)

Scenedesmus Subspicatus: 72hr IC50 >100 mg/l (OECD 201)

Mobility in Soil

No data available.

Bio-accumulative Potential

Not bioaccumulating.

Persistence and Degradability

Not readily biodegradable.

Other Adverse Effect

No data available.

Redux-823 Page 4 of 5

SECTION 13) DISPOSAL CONSIDERATIONS

Waste Disposal

Under RCRA it is the responsibility of the user of the product to determine at the time of disposal whether the product meets RCRA criteria for hazardous waste. Waste management should be in full compliance with federal, state and local laws. Empty Containers retain product residue which may exhibit hazards of material, therefore do not pressurize, cut, glaze, weld or use for any other purposes. Return drums to reclamation centers for proper cleaning and reuse.

SECTION 14) TRANSPORT INFORMATION

U.S. DOT Information

For all transportation accidents, call CHEMTREC at 800/424-9300. All spills and leaks of this material must be handled in accordance with local, state, and federal regulations.

DOT Shipping Designation:

Non-hazardous under 29-CFR 1910.1200. Water treatment compound

SECTION 15) REGULATORY INFORMATION

CAS	Chemical Name	% By Weight	Regulation List
No applicable CAS	No applicable chemical	-	-

SECTION 16) OTHER INFORMATION

Glossary

ACGIH- American Conference of Governmental Industrial Hygienists; ANSI- American National Standards Institute; Canadian TDGCanadian Transportation of Dangerous Goods; CAS- Chemical Abstract Service; Chemtrec- Chemical Transportation Emergency Center(US); CHIP- Chemical Hazard Information and Packaging; DSL- Domestic Substances List; EC- Equivalent Concentration; EH40 (UK)- HSE Guidance Note EH40 Occupational Exposure Limits; EPCRA- Emergency Planning and Community Right-To-Know Act; ESL Effects screening levels; HMIS- Hazardous Material Information Service; LC- Lethal Concentration; LD- Lethal Dose; NFPA- National Fire Protection Association; OEL- Occupational Exposure Limits; OSHA- Occupational Safety and Health Administration, US Department of Labor; PEL- Permissible Exposure Limit; SARA (Title III)- Superfund Amendments and Reauthorization Act; SARA 313- Superfund Amendments and Reauthorization Act, Section 313; SCBA- Self Contained Breathing Apparatus; STEL-Short Term Exposure Limit; TCEQ Texas Commission on Environmental Quality; TLV- Threshold Limit Value; TSCA- Toxic Substances Control Act Public Law 94-469; TWA Time Weighted Value; US DOT- US Department of Transportation; WHMIS- Workplace Hazardous Materials Information System.

Additional Information

Any concentration shown as a range is to protect confidentiality or is due to batch variation.

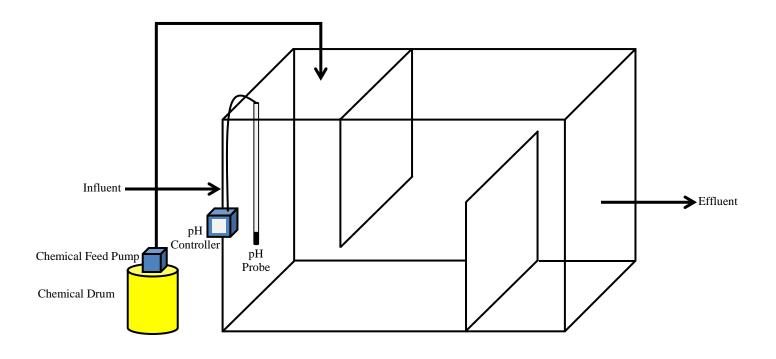
Version 1.0:

Revision Date: Jan 25, 2018 First Edition.

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Notes:

- 1.) Figure is not to scale.
- 2.) System layout can vary with site conditions.



89 Crawford Street Leominster, Massachusetts 01453

Tel: 774.450.7177 Fax: 888.835.0617 www.lrt-llc.net Configuration of pH Adjustment System



Lockwood Remediation Technologies, LLC



One Controller for the Broadest Range of Sensors.

Choose from 30 digital and analog sensor families for up to 17 di:erent parameters.

Maximum Versatility

The sc200 controller allows the use of digital and analog sensors, either alone or in combination, to provide compatibility with Hach's broad range of sensors, eliminating the need for dedicated, parameter-specific controllers.

Ease of Use and Confidence in Results

Large, high-resolution, transreflective display provides optimal viewing resolution in any lighting condition. Guided calibration procedures in 19 languages minimize complexity and reduce operator error. Password-protected SD card reader o:ers a simple solution for data download and transfer. Visual warning system provides critical alerts.

Wide Variety of Communication Options

Utilize two to five analog outputs to transmit primary and secondary values for each sensor, or integrate Hach sensors and analyzers into MODBUS RS232/RS485, Profibus® DP, and HART networks.



Password protected SD card reader offers a simple solution for data download and transfer, and sc200 and digital sensor configuration file duplication and backup.

Controller Comparison





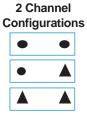


	Previous I	Models		
Features	sc100™ Controller	GLI53 Controller	sc200™ Controller	Benefits
Display	64 x 128 pixels 33 x 66 mm (1.3 x 2.6 in.)	64 x 128 pixels 33 x 66 mm (1.3 x 2.6 in.)	160 x 240 pixels 48 x 68 mm (1.89 x 2.67 in.) Transreflective	 Improved user interface— 50% bigger Easier to read in daylight and sunlight
Data Management	irDA Port/PDA Service Cable	N/A	SD Card Service Cable	 Simplifies data transfer Standardized accessories/ max compatibility
Sensor Inputs	2 Max Direct Digital Analog via External Gateway	2 Max Analog Depending on Parameter	2 Max Digital and/or Analog with Sensor Card	Simplifies analog sensor connectionsWorks with analog and digital sensors
Analog Inputs	N/A	N/A	1 Analog Input Signal Analog 4-20mA Card	 Enables non-sc analyzer monitoring Accepts mA signals from other analyzers for local display Consolidates analog mA signals to a digital output
4-20 mA Outputs	2 Standard	2 Standard	2 Standard Optional 3 Additional	Total of five (5) 4-20 mA outputs allows multiple mA outputs per sensor input
Digital Communication	MODBUS RS232/RS485 Profibus DP V1.0	HART	MODBUS RS232/RS485 Profibus DP V1.0 HART7.2	Unprecedented combination of sensor breadth and digital communication options

Choose from Hach's Broad Range of Digital and Analog Sensors					
Parameter	Sensor	Digital or Analog			
Ammonia	AMTAX™ sc, NH4D sc, AISE sc, AN-ISE sc	•			
Chlorine	CLF10 sc, CLT10 sc, 9184 sc	•			
Chlorine Dioxide	9185 sc	•			
Conductivity	GLI 3400 Contacting, GLI 3700 Inductive	A			
Dissolved Oxygen	LDO® Model 2, 5740 sc	•			
Dissolved Oxygen	5500	A			
Flow	U53, F53 Sensors	A			
Nitrate	NITRATAX™ sc, NO3D sc, NISE sc, AN-ISE sc	•			
Oil in Water	FP360 sc	•			
Organics	UVAS sc	•			
Ozone	9187 sc	•			
pH/ORP	pHD	•			
pH/ORP	pHD, pH Combination, LCP				
Phosphate	PHOSPHAX™ sc	•			
Sludge Level	SONATAX™sc	•			
Suspended Solids	SOLITAX™ sc, TSS sc	•			
Turbidity	1720E, FT660 sc, SS7 sc, ULTRATURB sc, SOLITAX sc, TSS sc	•			
Ultra Pure Conductivity	8310, 8311, 8312, 8315, 8316, 8317 Contacting	A			
Ultra Pure pH/ORP	8362	A			

● = Digital ▲ = Analog

Connect up to two of any of the sensors listed above, in any combination, to meet your application needs. The diagrams below demonstrate the potential configurations. Operation of analog sensors requires the controller to be equipped with the appropriate sensor module. Contact Hach Technical Support for help with selecting the appropriate module.



1 Channel
Configurations

Specifications*

Dimensions (H x W x

D)

5.7 in x 5.7 in x 7.1 in (144 mm x 144 mm x 181 mm)

Display

Weight

Graphic dot matrix LCD with LED backlighting, transreflective

Display Size

1.9 x 2.7 in. (48 mm x 68 mm)

Display Resolution

240 x 160 pixels 3.75 lbs. (1.70 kg)

Power Requirements

(Voltage)

Power Requirements

50/60 Hz

(Hz)

Operating

Temperature Range

Analog Outputs

-20 to 60 °C, 0 to 95% RH non-condensing

100 - 240 V AC, 24 V DC

Two (Five with optional expansion module) to isolated current

outputs, max 550 Ω , Accuracy: ± 0.1% of FS (20mA) at 25 °C, \pm 0.5% of FS over -20 °C to 60 °C

range

Operational Mode: measurement

or calculated value

Analog Output Functional Mode Linear, Logarithmic, Bi-linear, PID

Security Levels

Mounting **Configurations** 2 password-protected levels

Wall, pole, and panel mounting

Enclosure Rating Conduit Openings

Relay: Operational

Mode

1/2 in NPT Conduit

NEMA 4X/IP66

Primaryorsecondary

measurement, calculated value (dual channel only) or timer

Relay Functions

Scheduler (Timer), Alarm, Feeder Control, Event Control, Pulse Width Modulation, Frequency Control,

and Warning

Relays

Four electromechanical SPDT (Form C) contacts, 1200 W, 5 A

Communication MODBUS RS232/RS485,

PROFIBUS DPV1, or HART7.2

optional

Memory Backup

Electrical Certifications Flash memory

EMC

CE compliant for conducted and

radiated emissions:

- CISPR 11 (Class A limits)

- EMC Immunity EN 61326-1 (Industrial limits)

cETLus safety mark for:

- General Locations per ANSI/UL 61010-1 & CAN/CSA C22.2. No.

61010-1

Safety

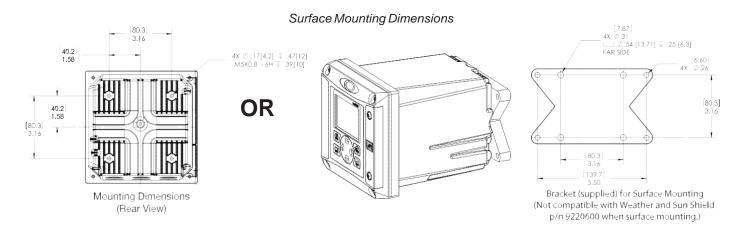
- Hazardous Location Class I, Division 2, Groups A,B,C & D (Zone 2, Group IIC) per FM 3600 / FM 3611 & CSA C22.2 No. 213 M1987 with approved options and appropriately rated Class I, Division 2 or Zone 2 sensors

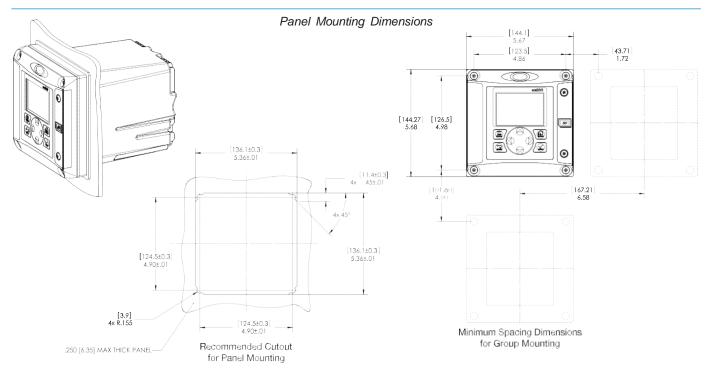
cULus safety mark

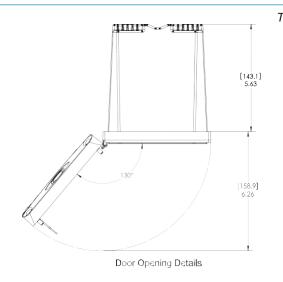
- General Locations per UL 61010-1 & CAN/CSA C22.2. No. 61010-1

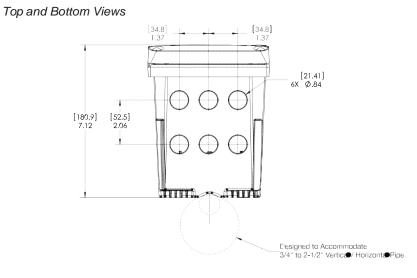
*Subject to change without notice.

Dimensions











3/4-inch Combination pH and ORP Sensor Kits





Use the Digital Gateway to make any Hach analog combination pH or ORP sensor compatible with the Hach sc1000 Controller.





Digital combination pH and ORP sensors are available in convertible, insertion, and sanitary mounting styles. Choose from rugged dome electrodes or "easy-to-clean" flat glass electrodes.

Features and Benefits

Low Price—High Performance

These combination sensors are designed for specialty applications for immersion or in-line mounting. The reference cell features a double-junction design for extended service life, and a built-in solution ground. The body is molded from chemically-resistant Ryton® or PVDF, and the reference junction is coaxial porous Teflon®. All sensors are rated 0 to 105°C up to 100 psig, and have integral 4.5 m (15 ft.) cables with tinned leads. The PC-series (for pH) and RC-series (for ORP) combination sensors are ideal for measuring mild and aggressive media.

Special Electrode Configurations

Sensors with rugged dome electrodes, "easy-to-clean" flat glass electrodes, and even HF (hydrofluoric acid) resistant glass electrodes are available for a wide variety of process solutions.

Temperature Compensation Element Option

The PC-series combination pH sensors are available with or without a Pt 1000 ohm RTD temperature element. The RC-series combination ORP sensors are supplied without a temperature element.

Versatile Mounting Styles

Sensors are available in three mounting styles—convertible, insertion, and sanitary. Please turn to page 3 for more information.

Full-Featured "Plug and Play" Hach sc Digital Controllers

There are no complicated wiring or set up procedures with any Hach sc controller. Just plug in any combination of Hach digital sensors and it's ready to use—it's "plug and play."

One or multiple sensors—The sc controller family allows you to receive data from up to eight Hach digital sensors in any combination using a single controller.

Communications—Multiple alarm/control schemes are available using the relays and PID control outputs. Available communications include analog 4-20 mA, digital MODBUS[®] (RS485 and RS232) or Profibus DP protocols. (Other digital protocols are available. Contact your Hach representative for details.)

Data logger—A built-in data logger collects measurement data, calibration, verification points, and alarm history.

Specifications*

Most pH applications fall in the 2.5-12.5 pH range. General purpose pH glass electrodes perform well in this range. Some industrial applications require accurate measurements and control at pH values below 2 or above 12. Consult Hach Technical Support for details on these applications.

Combination pH Sensors

Measuring Range

0 to 14 pH

Accuracy

Less than 0.1 pH under reference conditions

Temperature Range

0 to 105°C (32 to 221°F)

Flow Rate

0 to 2 m/s (0 to 6.6 ft./s); non-abrasive

Pressure Range

0 to 6.9 bar at 100°C (0 to 100 psig at 212°F)

Signal Transmission Distance

100 m (328 ft.) when used with the Hach Digital Gateway and a Hach sc Digital Controller.

1000 m (3280 ft.) when used with the Hach Digital Gateway, Termination Box, and a Hach sc Digital Controller.

Sensor Cable

Integral coaxial cable (plus two conductors for temperature compensator option); 4.5 m (15 ft.) long

Wetted Materials

Convertible style: Ryton® body (glass filled)

Insertion style: PVDF body (Kynar®)

Sanitary style: 316 stainless steel sleeved PVDF body

Common materials for all sensor styles include PTFE Teflon double junction, glass process electrode, and Viton® O-rings

Warranty

90 days

Combination ORP Sensors

Measuring Range

-2000 to +2000 millivolts

Accuracy

Limited to calibration solution accuracy (± 20 mV)

Temperature Range

0 to 105°C (32 to 221°F)

Flow Rate

0 to 2 m/s (0 to 6.6 ft./s); non-abrasive

Pressure Range

0 to 6.9 bar at 100°C (0 to 100 psig at 212°F)

Signal Transmission Distance

100 m (328 ft.) when used with the Hach Digital Gateway and a Hach sc Digital Controller.

1000 m (3280 ft.) when used with the Hach Digital Gateway, Termination Box, and a Hach sc Digital Controller.

Sensor Cable

Integral coaxial cable; 4.5 m (15 ft.) long; terminated with stripped and tinned wires

Wetted Materials

Convertible style: Ryton® body (glass filled)

Insertion style: PVDF body (Kynar®)

Common materials for all sensor styles include PTFE Teflon double junction, glass with platinum process electrode, and Viton® O-rings

Warranty

90 days

*Specifications subject to change without notice.

Ryton® is a registered trademark of Phillips 66 Co.; Viton® is a registered trademark of E.I. DuPont de Nemours + Co.; Kynar® is a registered trademark of Pennwalt Corp.

Engineering Specifications

- The pH sensor shall be available in convertible, insertion or sanitary styles. The ORP sensor shall be available in only convertible or insertion styles.
- 2. The convertible style sensor shall have a Ryton[®] body. The insertion style sensor shall have a PVDF body. The sanitary style sensor shall have a 316 stainless steel sleeved PVDF body. Common materials for all sensor styles shall include a PTFE Teflon[®] double junction, and Viton[®] O-rings. The pH sensor shall have a glass pH electrode. The ORP sensor shall have a platinum ORP electrode.
- The convertible style pH sensor shall be available with or without a built-in Pt 1000 ohm RTD temperature element. Insertion and sanitary style pH sensors shall have a built-in Pt 1000 ohm RTD temperature element. Convertible and insertion style ORP sensors shall not have a built-in temperature element.
- The sensor shall communicate via MODBUS[®] RS-485 to a Hach sc Digital Controller.
- The sensor shall be Hach Company Model PC sc or PC-series for pH measurement or Model PC sc or RC-series for ORP measurement.

Dimensions

Convertible Style Sensor

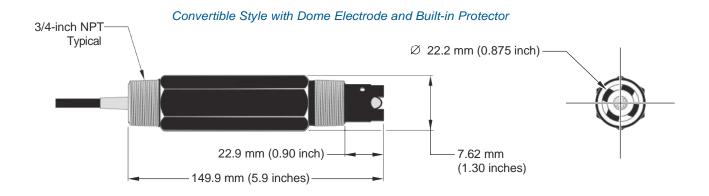
The convertible style sensor has a Ryton[®] body that features 3/4-inch NPT threads on both ends. The sensor can be directly mounted into a standard 3/4-inch pipe tee for flow-through mounting or fastened onto the end of a pipe for immersion mounting. The convertible style sensor enables inventory consolidation, thereby reducing associated costs. Mounting tees and immersion mounting hardware are offered in a variety of materials to suit application requirements.

Insertion Style Sensor

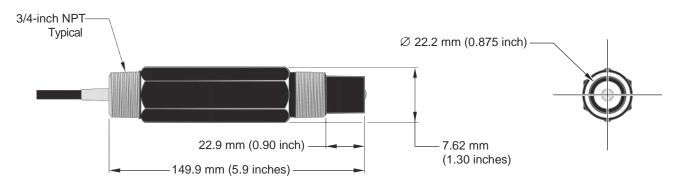
Insertion style sensors feature a longer, non-threaded PVDF body with two Viton® O-rings, providing a seal when used with the optional Hach insertion mount hardware assembly. This ball valve hardware enables sensor insertion and retraction from a pipe or vessel without having to stop the process flow.

Sanitary Style Sensor

The sanitary style sensor, offered for pH measurement, has a 316 stainless steel-sleeved PVDF body with a 2-inch flange. The sensor mates to a standard 2-inch Tri-Clover fitting. The optional Hach sanitary mounting hardware includes a standard 2-inch sanitary tee, sanitary clamp, and Viton[®] sanitary gasket.



Convertible Style with Flat Electrode





The Pulsatron Series A Plus offers manual function controls over stroke length and stroke rate as standard with the option to select external pace for automatic control.

Ten distinct models are available, having pressure capabilities to 250 PSIG (17 BAR) @ 12 GPO (1.9 lph), and flow capacities to 58 GPO (9.1 lph) @ 100 PSIG (7.0 BAR), with a standard turndown ratio of 100:1, and optional ratio of 1000:1. Metering performance is reproducible to within \pm 3% of maximum capacity.

Features

- Manual Control by on-line adjustable stroke rate and stroke length.
- Highly Reliable timing circuit.
- Circuit Protection against voltage and current upsets.
- Solenoid Protection by thermal overload with autoreset.
- Water Resistant, for outdoor and indoor applications.
- Internally Dampened To Reduce Noise.
- Guided Ball Check Valve Systems, to reduce back flow and enhance outstanding priming characteristics.
- Few Moving Parts and Wall Mountable.
- Safe & Easy Priming with durable leak-free bleed valve assembly (standard).
- Optional Control: External pace with auto/manual selection.

Controls



Manual Stroke Rate

Manual Stroke Length

External Pacing-Optional

External Pace With Stop-Optional (125 SPM only)

Controls Options						
F	Standard	Optional				
Feature	Configuration	Configuration ¹				
External Pacing		Auto / Manual Selection /				
External Pace w/ Stop		Auto / Manual Selection 2				
(125SPMonly)						
Manual Stroke Rate	10:1Ratio	100:1 Raio				
Manual Stroke Length	10:1 Ratio	10:1 Ratio				
Total Turndown Ratio	1001 Ratio	1000:1 Ratio				

Note 1:On S2,S3 & S4 sizes only.

Note 2:Not available on 1000:1turndown pumps.

Operating Benefits

- Reliable metering performance.
- Rated "hot" for continuous duty.
- High viscosity capability.
- Leak-free, sealless, liquid end.



Aftermarket

- KOPkits
- Gauges
- Dampeners
- Pressure Relief Valves
- Tanks
- Pre-Engineered Systems
- Process Controllers
 (PULSAblue, MicroVision)









Series A Plus Electronic Metering Pumps



Series A Plus

Specifications and Model Selection

	MODEL		LBC2	LB02	LBC3	LB03	LB04	LB64	LBC4	LBS2	LBS3	LBS4
Capacity		GPH	0.25	025	0.42	0.50	1.00	125	2.00	0.50	1.38	2.42
nominal		GPO	6	6	10	12	24	30	48	12	33	58
(max.)		LPH	0.9	0.9	1.6	1.9	3.8	4.7	7.6	1.9	5.2	9.14
Pressure ³ (max.)	GFPP,PVDF,316SS or PVC <;Ncode) wTFE Seats) PVC (V code) Vton or CSPE Seats IDegas Liquid End	PSIG (Bar)	250 (17) 150 (10)	150 (10)	250 (17)	150 (10)	100 (7)	100 (7)	50 (33)	250 (17) 150 (10)	150 (10)	100 (7)
Connections:		Tubina			114'IDX	318' OD			318'DX 112'OD	114	I'D X 318' OI)
		Pioina					1	14'FNPT				
Strokes/Minute		SPM		125					250			

Note 3: Pumps with rated pressure above 150 PSI will be de-rated to 150 PSI Max. when selecting certain valve options, see Price Book for details.

Engineering Data

Pump Head Materials Available: **GFPPL**

PVC **PVDF** 316 SS

Diaphragm: PTFE-faced CSPE-backed

Check Valves Materials Available:

Fittings Materials Available:

Seats/0-Rings: **PTFE**

> **CSPE** Viton

Balls: Ceramic

PTFE 316 SS

Alloy C **GFPPL**

PVC **PVDF**

Bleed Valve: Same as fitting and check valve

selected, except 316SS

hjection Valve & Foot Valve Assy: Same as fitting and check valve

selected

ClearPVC Tubing:

White PE

Important: Material Code - GFPPL=Glass-filled Polypropylene, PVC=Polyvinyl Chloride, PE=Polyethylene, PVDF=Polyvinylidene Fluoride, CSPE=Generic formulation of Hypalon, a registered trademark of E.I. DuPont Company. Viton is a registered trademark of E.I. DuPont Company. PVC wetted end recommended for sodium hypochlorite.

Engineering Data

Reproducibility: +/- 3% at maximum capady

Viscosity Max CPS: 1000CPS Stroke Frequency Max SPM: 125 / 250 by Model Stroke Frequency Turn-Down Ratio: 10:1/100:1 by Model

Stroke Length Turn-Down Ratio:

Power Input: 115 VAC/50-60 HZ/1 ph 230 VAC/50-60 HZ/1 ph

Average Current Draw:

@ 115 VAC; Amps: 0.6 Amps @ 230 VAC; Amps: 0.3 Amps 130 Watts Peak hout Power: 50 Watts Average Input Power @ Max SPM:

Custom Engineered Designs-Pre-Engineered Systems



Pre-Engineered Systems Pulsafeeder's Pre-Engineered Systems are designed to provide complete chemical feed solutions for all electronic metering applications. From stand alone simplex pH control applications to full-featured, redundant sodium hypochlorite disinfection metering, these rugged fabricated assemblies offer turnkey simplicity and industrial-grade durability. The UV-stabilized, high-grade HOPE frame offers maximum chemical compatibility and structural rigidity. Each system is factory assembled and hydrostatically tested prior to shipment.

Dimensions

Series A PLUS Dimensions (inches)							
Model No.	Α	В	С	D	Е	Weight	
LB02 IS2	5.0	9.6	9.5	6.5	8.2	10	
LBC2	5.0	9.9	9.5	6.5	8.5	10	
LBC3	5.0	9.9	9.5	6.5	8.5	10	
LB03 IS3	5.0	9.9	9.5	6.5	8.5	10	
LB0 \$ 4	5.0	9.9	9.5	6.5	8.5	10	
LB64	5.0	9.9	9.5	6.5	8.5	10	
LBC4	5.0	9.9	9.5	6.5	8.5	10	

NOTE: hches X 2.54 cm

BLEED VALVE	T T
	1AI



95-Gallon OverPack - 32" dia x 41.5", 1 each/package



Stock a SpillTech® OverPack with sorbents for emergency spill response, or use it as a salvage drum to ship damaged containers or hazardous waste.

- DOT-Approved for Salvage: All SpillTech® OverPacks are DOT-approved and X-rated for use as salvage drums. Helps companies conform to federal regulations when shipping damaged or leaking containers of hazardous materials, or absorbents contaminated with hazardous substances.
- Perfect for Spill Kits: Stores sorbent products (not included) for easy access as needed for spill control. Saves time when quick response is necessary.
- Sturdy Construction: 100% polyethylene OverPack resists chemicals, rust and corrosion for years of use. Integrated handles make them easy to lift, move or carry with standard material handling equipment. Twist-on, double-wall lid with closed-cell gasket provides sealed, secure closure to prevent leaks and protect contents from moisture, dirt and damage. Durable to withstand rough handling.
- Customized for You: We can customize a Spill Kit to your exact specifications, including the container, its contents and accessories, with no upcharge! Contact your local Distributor for details.

A950VER Specifications

Dimensions: ext. dia. 32" x 41.5" H

Shipping 31.75" W x 41.5" L x 31.75" H

Dimensions:

Sold as: 1 per package

Color: Yellow

Composition: Polyethylene

per Pallet: 3
Incinerable: No
Ship Class: 250

Metric Equivalent Specifications

Dimensions: ext. dia. 81.3cm x 105.4cm H

Shipping 80.6cm W x 105.4cm L x 80.6cm H

Dimensions:





A950VER Technical Information

Warnings & Restrictions:

There are no known warnings and restrictions for this product.

Regulations and Compliance:

49 CFR 173.3(c)(1) - If a container of hazardous waste is damaged or leaking, it can be placed in a compatible salvage drum that meets UN criteria for shipping

49 CFR 173.12(b)(2)(iv) - When labpacking, "Inner packagings...must be surrounded by a chemically compatible absorbent material in sufficient quantity to absorb the total liquid contents."

49 CFR 173.12(b) - A container used for labpacking must be "a UN 1A2 or UN 1B2 metal drum, a UN 1D plywood drum, a UN 1G fiber drum or a UN 1H2 plastic drum tested and marked at least for the Packing Group III performance level for liquids or solids."



Office: 774-450-7177 • Fax: 888-835-0617



Safety Data Sheet

According to U.S. Federal Register / Vol. 77, No. 58 / Monday, March 26, 2012 / Rules And Regulations and according to Canada's Hazardous Products Regulation, February 11, 2015.

Version: 4.0

Revision Date: 05/07/2018 Date of Issue: 05/31/2016

SECTION 1: IDENTIFICATION

Product Identifier

Product Form: Mixture

Product Name: Sulfuric Acid, 70-100%

Formula: H2-O4-S

Intended Use of the Product

Use Of The Substance/Mixture: Industrial use.

Name, Address, and Telephone of the Responsible Party

Manufacturer

CHEMTRADE LOGISTICS INC. 155 Gordon Baker Road

Suite 300

Toronto, Ontario M2H 3N5
For SDS Info: (416) 496-5856
www.chemtradelogistics.com
Emergency Telephone Number

Emergency Number : Canada: CANUTEC +1-613-996-6666 / US: CHEMTREC +1-800-424-9300

INTERNATIONAL: +1-703-741-5970 Chemtrade Emergency Contact: (866) 416-4404

For Chemical Emergency, Spill, Leak, Fire, Exposure, or Accident, call CHEMTREC - Day or Night

SECTION 2: HAZARDS IDENTIFICATION

Classification of the Substance or Mixture

GHS Classification

 Met. Corr. 1
 H290

 Skin Corr. 1A
 H314

 Eye Dam. 1
 H318

 Carc. 1A
 H350

 Aquatic Acute 3
 H402

Full text of hazard classes and H-statements: see section 16

Label Elements

GHS Labeling

Hazard Pictograms





Signal Word : Danger

Hazard Statements : H290 - May be corrosive to metals.

H314 - Causes severe skin burns and eye damage.

H318 - Causes serious eye damage. H350 - May cause cancer (Inhalation).

H402 - Harmful to aquatic life.

Precautionary Statements : P201 - Obtain special instructions before use.

P202 - Do not handle until all safety precautions have been read and understood.

P234 - Keep only in original container.

P260 - Do not breathe vapors, mist, or spray.

P264 - Wash hands, forearms, and other exposed areas thoroughly after handling.

P273 - Avoid release to the environment.

P280 - Wear protective gloves, protective clothing, and eye protection. P301+P330+P331 - IF SWALLOWED: Rinse mouth. Do NOT induce vomiting.

05/07/2018 EN (English US) SDS#: CHE-1010S 1/9

Safety Data Sheet

According to U.S. Federal Register / Vol. 77, No. 58 / Monday, March 26, 2012 / Rules And Regulations and according to Canada's Hazardous Products Regulation, February 11, 2015.

P303+P361+P353 - IF ON SKIN (or hair): Take off immediately all contaminated clothing. Rinse skin with water.

P304+P340 - IF INHALED: Remove person to fresh air and keep comfortable for breathing.

P305+P351+P338 - IF IN EYES: Rinse cautiously with water for several minutes. Remove contact lenses, if present and easy to do. Continue rinsing.

P308+P313 - If exposed or concerned: Get medical advice/attention.

P310 - Immediately call a POISON CENTER or doctor.

P321 - Specific treatment (see section 4 on this SDS).

P363 - Wash contaminated clothing before reuse.

P390 - Absorb spillage to prevent material damage.

P405 - Store locked up.

P406 - Store in corrosive resistant container with a resistant inner liner.

P501 - Dispose of contents/container in accordance with local, regional, national,

territorial, provincial, and international regulations.

Other Hazards

Exposure may aggravate pre-existing eye, skin, or respiratory conditions.

Unknown acute toxicity

No data available

SECTION 3: COMPOSITION/INFORMATION ON INGREDIENTS

Mixture

Name	Product Identifier	%*	GHS Ingredient Classification
Sulfuric acid**	(CAS-No.) 7664-93-9	70 - 100	Met. Corr. 1, H290
			Skin Corr. 1A, H314
			Eye Dam. 1, H318
			Carc. 1A, H350
			Aquatic Acute 3, H402
Water	(CAS-No.) 7732-18-5	0.1 - 30	Not classified

Full text of H-phrases: see section 16

SECTION 4: FIRST AID MEASURES

Description of First-aid Measures

General: Never give anything by mouth to an unconscious person. If you feel unwell, seek medical advice (show the label where possible).

Inhalation: When symptoms occur: go into open air and ventilate suspected area. Obtain medical attention if breathing difficulty persists.

Skin Contact: Remove contaminated clothing. Immediately flush skin with plenty of water for at least 30 minutes. Get immediate medical advice/attention. Wash contaminated clothing before reuse.

Eye Contact: Rinse cautiously with water for at least 30 minutes. Remove contact lenses, if present and easy to do. Continue rinsing. Get immediate medical advice/attention.

Ingestion: Rinse mouth. Do NOT induce vomiting. Obtain medical attention.

Most Important Symptoms and Effects Both Acute and Delayed

General: Corrosive to eyes, respiratory system and skin. May cause cancer.

Inhalation: May be corrosive to the respiratory tract.

Skin Contact: Causes severe irritation which will progress to chemical burns. **Eye Contact:** Causes permanent damage to the cornea, iris, or conjunctiva.

Ingestion: May cause burns or irritation of the linings of the mouth, throat, and gastrointestinal tract.

05/07/2018 EN (English US) SDS#: CHE-1010S 2/9

^{*}Percentages are listed in weight by weight percentage (w/w%) for liquid and solid ingredients. Gas ingredients are listed in volume by volume percentage (v/v%).

^{**}Strong inorganic acid aerosols/mists containing this substance are carcinogenic to humans via inhalation. Under normal conditions of use this route of exposure is not expected.

Safety Data Sheet

According to U.S. Federal Register / Vol. 77, No. 58 / Monday, March 26, 2012 / Rules And Regulations and according to Canada's Hazardous Products Regulation, February 11, 2015.

Chronic Symptoms: Strong inorganic acid mists containing sulfuric acid are carcinogenic to humans. Prolonged inhalation of fumes or mists may cause erosion of the teeth.

<u>Indication of Any Immediate Medical Attention and Special Treatment Needed</u>

If exposed or concerned, get medical advice and attention. If medical advice is needed, have product container or label at hand.

SECTION 5: FIRE FIGHTING MEASURES

Extinguishing Media

Suitable Extinguishing Media: Foam, carbon dioxide, dry chemical.

Unsuitable Extinguishing Media: Do not use water. Do not get water inside containers. Do not apply water stream directly at source of leak

Special Hazards Arising From the Substance or Mixture

Fire Hazard: Not flammable.

Explosion Hazard: Product is not explosive.

Reactivity: May be corrosive to metals. Contact with metals may evolve flammable hydrogen gas. May react exothermically with water releasing heat. Adding an acid to a base or base to an acid may cause a violent reaction. This product may act as an oxidizer.

Advice for Firefighters

Precautionary Measures Fire: Exercise caution when fighting any chemical fire.

Firefighting Instructions: Use water spray or fog for cooling exposed containers.

Protection During Firefighting: Do not enter fire area without proper protective equipment, including respiratory protection.

Hazardous Combustion Products: Toxic fumes are released.

Other Information: Do not allow run-off from fire fighting to enter drains or water courses.

Reference to Other Sections

Refer to Section 9 for flammability properties.

SECTION 6: ACCIDENTAL RELEASE MEASURES

Personal Precautions, Protective Equipment and Emergency Procedures

General Measures: Do not get in eyes, on skin, or on clothing. Do not breathe vapor, mist or spray. Do not handle until all safety precautions have been read and understood.

For Non-Emergency Personnel

Protective Equipment: Use appropriate personal protective equipment (PPE).

Emergency Procedures: Evacuate unnecessary personnel.

For Emergency Personnel

Protective Equipment: Equip cleanup crew with proper protection.

Emergency Procedures: Upon arrival at the scene, a first responder is expected to recognize the presence of dangerous goods, protect oneself and the public, secure the area, and call for the assistance of trained personnel as soon as conditions permit. Ventilate area.

Environmental Precautions

Prevent entry to sewers and public waters. Avoid release to the environment.

Methods and Materials for Containment and Cleaning Up

For Containment: Contain any spills with dikes or absorbents to prevent migration and entry into sewers or streams. As an immediate precautionary measure, isolate spill or leak area in all directions.

Methods for Cleaning Up: Clean up spills immediately and dispose of waste safely. Absorb spillage to prevent material damage. Cautiously neutralize spilled liquid. Transfer spilled material to a suitable container for disposal. Contact competent authorities after a spill.

Reference to Other Sections

See Section 8 for exposure controls and personal protection and Section 13 for disposal considerations.

SECTION 7: HANDLING AND STORAGE

Precautions for Safe Handling

Wash hands and other exposed areas with mild soap and water before eating, drinking or smoking and when leaving work. Handle empty containers with care because they may still present a hazard. Do not get in eyes, on skin, or on clothing. Do not breathe vapors, mist, spray. Obtain special instructions before use. Do not handle until all safety precautions have been read and understood.

Additional Hazards When Processed: May be corrosive to metals. May release corrosive vapors. NEVER pour water into this substance; when dissolving or diluting always add it slowly to the water.

05/07/2018 EN (English US) SDS#: CHE-1010S 3/9

Safety Data Sheet

According to U.S. Federal Register / Vol. 77, No. 58 / Monday, March 26, 2012 / Rules And Regulations and according to Canada's Hazardous Products Regulation, February 11, 2015.

Hygiene Measures: Handle in accordance with good industrial hygiene and safety procedures.

Conditions for Safe Storage, Including Any Incompatibilities

Technical Measures: Comply with applicable regulations.

Storage Conditions: Keep container closed when not in use. Store in a dry, cool place. Keep/Store away from extremely high or low

temperatures and incompatible materials. Store in original container or corrosive resistant and/or lined container.

Incompatible Materials: Combustible materials. Reducing agents. Strong oxidizers. Strong bases. Metals. Water.

Specific End Use(s)

Industrial use.

SECTION 8: EXPOSURE CONTROLS/PERSONAL PROTECTION

Control Parameters

For substances listed in section 3 that are not listed here, there are no established Exposure limits from the manufacturer, supplier, importer, or the appropriate advisory agency including: ACGIH (TLV), AIHA (WEEL), NIOSH (REL), OSHA (PEL), Canadian provincial governments, or the Mexican government.

Sulfuric acid (7664-93-9)	-	
Mexico	OEL TWA (mg/m³)	1 mg/m³
USA ACGIH	ACGIH TWA (mg/m³)	0.2 mg/m³ (thoracic particulate matter)
USA ACGIH	ACGIH chemical category	Suspected Human Carcinogen contained in strong
		inorganic acid mists
USA OSHA	OSHA PEL (TWA) (mg/m³)	1 mg/m³
USA NIOSH	NIOSH REL (TWA) (mg/m³)	1 mg/m³
USA IDLH	US IDLH (mg/m³)	15 mg/m³
Alberta	OEL STEL (mg/m³)	3 mg/m³
Alberta	OEL TWA (mg/m³)	1 mg/m³
British Columbia	OEL TWA (mg/m³)	0.2 mg/m³ (Thoracic, contained in strong inorganic acid
		mists)
Manitoba	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic particulate matter)
New Brunswick	OEL STEL (mg/m³)	3 mg/m³
New Brunswick	OEL TWA (mg/m³)	1 mg/m³
Newfoundland & Labrador	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic particulate matter)
Nova Scotia	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic particulate matter)
Nunavut	OEL STEL (mg/m³)	0.6 mg/m³ (thoracic fraction)
Nunavut	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic fraction)
Northwest Territories	OEL STEL (mg/m³)	0.6 mg/m³ (thoracic fraction, strong acid mists only)
Northwest Territories	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic fraction, strong acid mists only)
Ontario	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic)
Prince Edward Island	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic particulate matter)
Québec	VECD (mg/m³)	3 mg/m³
Québec	VEMP (mg/m³)	1 mg/m³
Saskatchewan	OEL STEL (mg/m³)	0.6 mg/m³ (thoracic fraction)
Saskatchewan	OEL TWA (mg/m³)	0.2 mg/m³ (thoracic fraction)
Yukon	OEL STEL (mg/m³)	1 mg/m³
Yukon	OEL TWA (mg/m³)	1 mg/m³

Exposure Controls

Appropriate Engineering Controls: Emergency eye wash fountains and safety showers should be available in the immediate vicinity of any potential exposure. Ensure adequate ventilation, especially in confined areas. Ensure all national/local regulations are observed.

Personal Protective Equipment: Gloves. Protective clothing. Protective goggles. Face shield. Insufficient ventilation: wear respiratory protection.











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Materials for Protective Clothing: Acid-resistant clothing.

Hand Protection: Wear protective gloves.

Eye Protection: Chemical safety goggles and face shield. **Skin and Body Protection:** Wear suitable protective clothing.

Respiratory Protection: If exposure limits are exceeded or irritation is experienced, approved respiratory protection should be worn. In case of inadequate ventilation, oxygen deficient atmosphere, or where exposure levels are not known wear approved respiratory protection.

Other Information: When using, do not eat, drink or smoke.

SECTION 9: PHYSICAL AND CHEMICAL PROPERTIES

Information on Basic Physical and Chemical Properties

Physical State : Liquid

Appearance : Clear, Colorless to Amber, Oily

Odor Threshold : Pungent : Not available

pH : 0.3

Evaporation Rate Not available **Melting Point** 10.56 °C (51.01 °F) **Freezing Point** Not available **Boiling Point** 290 °C (554 °F) **Flash Point** Not applicable **Auto-ignition Temperature** Not applicable **Decomposition Temperature** Not available Flammability (solid, gas) Not applicable **Lower Flammable Limit** Not applicable **Upper Flammable Limit** Not applicable

Vapor Pressure : 0.00027 - 0.16 kPa at 25 °C (77 °F)

Relative Vapor Density at 20°C : 3.4 (air = 1)
Relative Density : Not available
Specific Gravity : 1.84 g/l

Solubility : Water: Miscible
Partition Coefficient: N-Octanol/Water : Not available
Viscosity : Not available

SECTION 10: STABILITY AND REACTIVITY

Reactivity: May be corrosive to metals. Contact with metals may evolve flammable hydrogen gas. May react exothermically with water releasing heat. Adding an acid to a base or base to an acid may cause a violent reaction. This product may act as an oxidizer.

Chemical Stability: Stable under recommended handling and storage conditions (see section 7).

<u>Possibility of Hazardous Reactions</u>: Hazardous polymerization will not occur.

Conditions to Avoid: Extremely high or low temperatures and incompatible materials.

<u>Incompatible Materials:</u> Combustible materials. Reducing agents. Strong bases. Strong oxidizers. Metals. Water.

Hazardous Decomposition Products: Thermal decomposition generates: Corrosive vapors.

SECTION 11: TOXICOLOGICAL INFORMATION

Information on Toxicological Effects - Product

Acute Toxicity (Oral): Not classified
Acute Toxicity (Dermal): Not classified
Acute Toxicity (Inhalation): Not classified

LD50 and LC50 Data: Not available

Skin Corrosion/Irritation: Causes severe skin burns and eye damage.

pH: 0.3

Eye Damage/Irritation: Causes serious eye damage.

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pH: 0.3

Respiratory or Skin Sensitization: Not classified

Germ Cell Mutagenicity: Not classified

Carcinogenicity: May cause cancer (Inhalation).

Specific Target Organ Toxicity (Repeated Exposure): Not classified

Reproductive Toxicity: Not classified

Specific Target Organ Toxicity (Single Exposure): Not classified

Aspiration Hazard: Not classified

Symptoms/Effects After Inhalation: May be corrosive to the respiratory tract.

Symptoms/Effects After Skin Contact: Causes severe irritation which will progress to chemical burns. **Symptoms/Effects After Eye Contact:** Causes permanent damage to the cornea, iris, or conjunctiva.

Symptoms/Effects After Ingestion: May cause burns or irritation of the linings of the mouth, throat, and gastrointestinal tract.

Chronic Symptoms: Strong inorganic acid mists containing sulfuric acid are carcinogenic to humans. Prolonged inhalation of fumes or mists may cause erosion of the teeth.

<u>Information on Toxicological Effects - Ingredient(s)</u>

LD50 and LC50 Data:

Water (7732-18-5)				
LD50 Oral Rat	> 90000 mg/kg			
Sulfuric acid (7664-93-9)				
LD50 Oral Rat	2140 mg/kg			
C50 Inhalation Rat 510 mg/m³ (Exposure time: 2 h)				
Sulfuric acid (7664-93-9)				
IARC Group	1			
OSHA Hazard Communication Carcinogen List	In OSHA Hazard Communication Carcinogen list.			
Strong inorganic acid mists containing sulfuric acid				
National Toxicology Program (NTP) Status	Known Human Carcinogens.			

SECTION 12: ECOLOGICAL INFORMATION

Toxicity

Ecology - General: Harmful to aquatic life.

Sulfuric acid (7664-93-9)	
LC50 Fish 1	500 mg/l (Exposure time: 96 h - Species: Brachydanio rerio [static])
LC50 Fish 2	42 mg/l (Exposure time: 96 h - Species: Gambusia affinis [static])

Persistence and Degradability

Sulfuric Acid, 70-100%	
Persistence and Degradability	Not established.

Bioaccumulative Potential

Sulfuric Acid, 70-100%	
Bioaccumulative Potential	Not established.
Sulfuric acid (7664-93-9)	
BCF Fish 1	(no bioaccumulation)

Mobility in Soil Not available

Other Adverse Effects

Other Information: Avoid release to the environment.

SECTION 13: DISPOSAL CONSIDERATIONS

Waste Disposal Recommendations: Dispose of contents/container in accordance with local, regional, national, territorial, provincial, and international regulations.

Additional Information: Container may remain hazardous when empty. Continue to observe all precautions.

Ecology - Waste Materials: Avoid release to the environment. This material is hazardous to the aquatic environment. Keep out of sewers and waterways.

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SECTION 14: TRANSPORT INFORMATION

The shipping description(s) stated herein were prepared in accordance with certain assumptions at the time the SDS was authored, and can vary based on a number of variables that may or may not have been known at the time the SDS was issued.

TRANSPORTATION	DOT	TDG	IMDG	IATA
CLASSIFICATION				
Identification Number	UN1830	UN1830	UN1830	UN1830
Proper Shipping Name	SULFURIC ACID	SULFURIC ACID	SULPHURIC ACID	SULPHURIC ACID
Transport Hazard	8	8	8	8
Class(es)				
	CORROSIVE			0
Packing Group	II	II	II	II
Environmental Hazards	Marine Pollutant : No	Marine Pollutant : No	Marine Pollutant : No	Marine Pollutant: N/A
Emergency Response	ERG Number: 137	ERAP Index: 3 000	EMS: F-A, S-B	ERG code (IATA): 8L
Additional Information	Not applicable	Not applicable	Not applicable	Not applicable

SECTION 15: REGULATORY INFORMATION

US Federal Regulations

Chemical Name (CAS No.)	CERCLA RQ	EPCRA 304 RQ	SARA 302 TPQ	SARA 313
Sulfuric acid (7664-93-9)	1000 lb	1000 lb	1000 lb	Yes

SARA 311/312

Sulfuric Acid, 70-100%

Immediate (acute) health hazard. Delayed (chronic) health hazard. Reactive hazard

US TSCA Flags Not present

US State Regulations

California Proposition 65

Chemical Name (CAS No.)	Carcinogenicity	Developmental Toxicity	Female Reproductive Toxicity	Male Reproductive Toxicity
Sulfuric acid (7664-93-9)	Yes	No	No	No
Strong inorganic acid mists containing sulfuric acid	Yes	No	No	No

State Right-To-Know Lists

Sulfuric acid (7664-93-9)

- U.S. Massachusetts Right To Know List Yes
- U.S. New Jersey Right to Know Hazardous Substance List Yes
- U.S. Pennsylvania RTK (Right to Know) Environmental Hazard List Yes
- U.S. Pennsylvania RTK (Right to Know) Special Hazardous Substances No
- U.S. Pennsylvania RTK (Right to Know) List Yes

Canadian Regulations

Sulfuric acid (7664-93-9)

Listed on the Canadian DSL (Domestic Substances List)

Not listed on the Canadian NDSL (Non-Domestic Substances List)

International Inventories/Lists

Chemical Name (CAS No.)	Australia	Turkey	Korea	EU	EU	EU	EU	Mexico
	AICS	CICR	ECL	EINECS	ELINCS	SVHC	NLP	INSQ
Sulfuric acid (7664-93-9)	Yes	No	Yes	Yes	No	No	No	No

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Chemical Name (CAS No.)	China IECSC	Japan ENCS	Japan ISHL	Japan PDSCL	Japan PRTR	Philippines PICCS	New Zealand NZIOC	US TSCA
Sulfuric acid (7664-93-9)	Yes	Yes	No	Yes	No	Yes	Yes	Yes

SECTION 16: OTHER INFORMATION, INCLUDING DATE OF PREPARATION OR LAST REVISION

Date of Preparation or Latest Revision : 05/07/2018

Revision Summary

Section	Change	Date Changed
16	Data modified	05/07/2018

Other Information

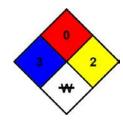
: This document has been prepared in accordance with the SDS requirements of the OSHA Hazard Communication Standard 29 CFR 1910.1200 and Canada's Hazardous Products Regulations (HPR).

GHS Full Text Phrases:

Aquatic Acute 3	Hazardous to the aquatic environment - Acute Hazard Category 3
Carc. 1A	Carcinogenicity Category 1A
Eye Dam. 1	Serious eye damage/eye irritation Category 1
Met. Corr. 1	Corrosive to metals Category 1
Skin Corr. 1A	Skin corrosion/irritation Category 1A
H290	May be corrosive to metals
H314	Causes severe skin burns and eye damage
H318	Causes serious eye damage
H350	May cause cancer
H402	Harmful to aquatic life

NFPA 704

NFPA Health Hazard : 3
NFPA Fire Hazard : 0
NFPA Reactivity Hazard : 2
NFPA Specific Hazards : W



HMIS Rating

Health : 3 Flammability : 0 Physical : 2

PPE See Section 8

ENCS - Japanese Existing and New Chemical Substances Inventory

Abbreviations and Acronyms

AICS – Australian Inventory of Chemical Substances LC50 - Median Lethal Concentration

ACGIH – American Conference of Governmental Industrial Hygienists LD50 - Median Lethal Dose

AIHA – American Industrial Hygiene Association

LOAEL - Lowest Observed Adverse Effect Level

ATE - Acute Toxicity Estimate

LOEC - Lowest-observed-effect Concentration

BEI - Biological Exposure Indices (BEI)

LOEC - Lowest-observed-effect Concentration

Log Pow - Octanol/water Partition Coefficient

NFPA 704 – National Fire Protection Association - Standard System for the

CAS No. - Chemical Abstracts Service number Identification of the Hazards of Materials for Emergency Response

CERCLA RQ - Comprehensive Environmental Response, Compensation, and
Liability Act - Reportable Quantity

NIOSH - National Institute for Occupational Safety and Health
NLP - Europe No Longer Polymers List

CICR - Turkish Inventory and Control of Chemicals

DOT – 49 CFR – US Department of Transportation – Code of Federal

Regulations Title 49 – Transportation.

NOAEL - No-Observed Adverse Effect Level

NOEC - No-Observed Effect Concentration

NZIOC - New Zealand Inventory of Chemicals

EC50 - Median effective concentration OEL - Occupational Exposure Limits

ECL - Korea Existing Chemicals List

OSHA – Occupational Safety and Health Administration

EINECS - European Inventory of Existing Commercial Chemical Substances

PEL - Permissible Exposure Limits

ELINCS - European List of Notified Chemical Substances

EmS - IMDG Emergency Schedule Fire & Spillage

PICCS - Philippine Inventory of Chemicals and Chemical Substances

PDSCL - Japan Poisonous and Deleterious Substances Control Law

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PPE - Personal Protective Equipment

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EPA – Environmental Protection Agency

EPCRA 304 RQ – EPCRA 304 Extremely Hazardous Substance Emergency Planning and Community Right-to-Know-Act – Reportable Quantity ERAP Index – Emergency Response Assistance Plan Quantity Limit

ErC50 - EC50 in Terms of Reduction Growth Rate

ERG code (IATA) - Emergency Response Drill Code as found in the International

Civil Aviation Organization (ICAO)

ERG No. - Emergency Response Guide Number HCCL - Hazard Communication Carcinogen List HMIS – Hazardous Materials Information System IARC - International Agency for Research on Cancer

IATA - International Air Transport Association – Dangerous Goods Regulations

IDLH - Immediately Dangerous to Life or Health

IECSC - Inventory of Existing Chemical Substances Produced or Imported in

China

IMDG - International Maritime Dangerous Goods Code INSQ - Mexican National Inventory of Chemical Substances

ISHL - Japan Industrial Safety and Health Law

PRTR - Japan Pollutant Release and Transfer Register

REL - Recommended Exposure Limit

SADT - Self Accelerating Decomposition Temperature SARA - Superfund Amendments and Reauthorization Act

SARA 302 - Section 302, 40 CFR Part 355

SARA 311/312 - Sections 311 and 312, 40 CFR Part 370 Hazard Categories

SARA 313 - Section 313, 40 CFR Part 372 SRCL - Specifically Regulated Carcinogen List

STEL - Short Term Exposure Limit

SVHC – European Candidate List of Substance of Very High Concern TDG – Transport Canada Transport of Dangerous Goods Regulations

TLM - Median Tolerance Limit TLV - Threshold Limit Value TPQ - Threshold Planning Quantity

TSCA - United StatesToxic Substances Control Act

TWA - Time Weighted Average

WEEL - Workplace Environmental Exposure Levels

Handle product with due care and avoid unnecessary contact. This information is supplied under U.S. OSHA'S "Right to Know" (29 CFR 1910.1200) and Canada's WHMIS regulations. Although certain hazards are described herein, we cannot guarantee these are the only hazards that exist. The information contained herein is based on data available to us and is believed to be true and accurate but it is not offered as a product specification. No warranty, expressed or implied, regarding the accuracy of this data, the hazards connected with the use of the product, or the results to be obtained from the use thereof, is made and Chemtrade and its affiliates assume no responsibility. Chemtrade is a member of the CIAC (Chemistry Industry Association of Canada) and adheres to the codes and principles of Responsible Care™.



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