Stantec Consulting Services, Inc. 400 Crown Colony Drive, Suite 200, Quincy, MA 02169



February 21, 2020 File: 195601892

Attention: US Environmental Protection Agency
Office of Ecosystem Protection

5 Post Office Square – Suite 100 (OEP06-01)

Boston, Massachusetts 02109

Attn: Shelley Puleo Shauna Little

EPA / OEP RGP Applications Coordinator

Reference: Remediation General Permit (RGP)

30 Penniman Road Allston, Massachusetts

On behalf of our client, 30 Penn, LLC, Stantec Consulting Services, Inc. has prepared this submission for a National Pollutant Discharge Elimination System (NPDES) Remediation General Permit (RGP) to facilitate off-site discharge of dewatering effluent generated during construction activities beneath 30 Penniman Road in Allston, Massachusetts (the Site, see Figure 1). The information presented herein has been prepared to follow requirements of the 2017 US Environmental Protection Agency (EPA) NPDES RGP. A copy of the completed Notice of Intent (NOI) form is enclosed as Appendix A.

Since the Site is a listed Massachusetts of Department of Environmental Protection (MassDEP) Massachusetts Contingency Plan (MCP) Disposal Site (discussed below), notification to the State is not required.

SITE HISTORY AND EXISTING CONDITIONS

The parcel associated with the Site consists of approximately 11,639 square feet of land with frontage along Penniman Road on the west side. The Site was part of a commercial property developed before 1925 by the Albany Carpet Cleaning Company. This company occupied the Site until around 1975 when the usage changed to commercial activities mainly associated with artist tenant rentals. The Site currently consists of a construction site which is being developed into a multi-story residential / commercial building. The Site associated with this NPDES RGP is delineated within the extent of the area of the parcel (see Figure 2).

The Albany Carpet Cleaning Company occupied 18 to 30 Penniman Road and was a carpet and upholstery cleaning business. A release at the former Albany Carpet Cleaning Company from former fuel oil USTs located in a courtyard adjacent to the north of the Site was managed under MassDEP Release Tracking Number (RTN) 3-16868. The majority of impacted soil was excavated from the accessible portions of the courtyard in early response actions. Some oil was not accessible since it had migrated as non-aqueous phase liquid (NAPL) beneath the former Site building and a former abutting building to the north. The abutting parcel to the north was redeveloped and underwent extensive excavation up to 10 feet below grade. The excavation extended up to the northern side of the Site. Subsequent assessment indicated that

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Reference: Remediation General Permit (RGP) 30 Penniman Road Allston, Massachusetts

the excavation at the abutting parcel eliminated the remaining petroleum NAPL beneath the former Site building.

A second DEP RTN 3-27600 was linked to the Albany Carpet Cleaning Company's Primary RTN 3-16868 for lead detected in soil samples collected beneath the former building. Similar lead impacts were observed within the excavated area of the abutting property. Subsequent evaluation indicated the presence of lead was likely the result of the placement of historic fill and not associated with a release.

Based upon the historic fill determination for the lead, and the elimination of the fuel oil and the waste oil NAPLs, the RTNs were closed with the submittal to MassDEP of a Permanent Solution Statement with No Conditions dated June 13, 2017.

NEW RELEASE NOTIFICATION

As part of soil characterization activities, Stantec collected soil samples across the Site on January 9, 2020. The samples were submitted to Phoenix Environmental Laboratory located in Manchester, Connecticut for waste characterization analyses. The analytical results indicate that mercury and zinc were detected is soil at concentrations that exceed the applicable MassDEP RCS-1 reportable concentrations. The impacts were limited to a surface fill layer. Accordingly, on February 3, 2020, 30 Penn, LLC notified the MassDEP and the RTN 3-36120 was issued. TCLP results for mercury were below the laboratory method detection limits (set below the EPA threshold) indicating the mercury is not likely to leach. This mercury and zincimpacted fill material is scheduled for excavation and off-Site disposal prior to the start of dewatering activities.

PROPOSED CONSTRUCTION

The proposed project consists of excavating the footprint of the parcel for construction of footings and an elevator shaft for the proposed building. The excavation will extend up to 20 feet below the current grade (approximate EL = 30 feet above NGVD). Depth to groundwater ranges between 6 and 11 feet below grade indicating excavation activities will need to be conducted below the water table.

CURRENT GROUNDWATER QUALITY DATA

To evaluate groundwater quality at the Site, two observation wells (C4 and D1 on Figure 2) were installed via the direct push drilling methodology. Groundwater samples were collected from the wells on January 10, 2020 and submitted to Phoenix Environmental Laboratory for analysis of RGP parameters. pH and temperature readings were collected in the field.

Results of the analyses indicated cis-1,2-dichloroethene, tetrachloroethylene, trichloroethylene, and vinyl chloride were detected at concentrations that exceed the applicable DEP Method 1 RCGW-2 reportable concentrations (see Table 1). These contaminants are associated with a nearby disposal site and are not due to releases at the Site so did not warrant a new notification to DEP. The nearby disposal site is known as the former Sunshine Laundry site and it is currently an active Massachusetts Contingency Plan (MCP) site under DEP RTN 3-0506. Copies of the laboratory reports are included in Appendix B. All other detected analytes were below the RCGW-2 reportable concentrations.

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Reference: Remediation General Permit (RGP) 30 Penniman Road Allston, Massachusetts

RECEIVING WATERS SAMPLING AND DILUTION FACTOR

On February 14, 2020, Stantec collected one surface water sample designated CHAR-SW from upstream of the proposed SDO037 outfall location into the Charles River. The sample was submitted to Phoenix for total metals, ammonia and hardness. pH and temperature readings were collected in the field. The analytical results indicate that ammonia (as nitrogen) was detected at 0.14 mg/L, hardness was 74.4 mg/L, and iron and zinc were detected in the sample at low concentrations (see Table 2). A copy of the laboratory report is included in Appendix B.

The seven-day-ten-year flow (7Q10) of the receiving water was established using the U.S. Geological Survey (USGS) StreamStats program and confirmed by Massachusetts Department of Environmental Protection (MassDEP) on February 12, 2020. We have additionally confirmed with the MassDEP that the dilution factor for the receiving waters is 22.81. The StreamStats Report, Dilution Factor calculations, and confirmation from MassDEP are included in Appendix C.

EFFLUENT CRITERIA DOCUMENTATION

Groundwater and Receiving Water data were input into the MALimitsBook calculation spreadsheet provided by EPA. Per a communication with Shauna Little at EPA, the "FreshwaterResults" tab of the downloadable spreadsheet is being reviewed by EPA and not available to calculate the effluent criteria for the Site. Ms. Little indicated these values will be provided by EPA during their review of the NOI. A copy of the "EnterData" tab of the spreadsheet is included in Appendix C.

DEWATERING SYSTEM AND OFF-SITE DISCHARGE

During construction activities, it will be necessary to perform temporary dewatering to control surface water runoff from groundwater seepage to enable construction in-the-dry. Construction and construction dewatering activities are currently anticipated to be required for a period of over 6 months. On average, we estimate effluent discharge rates of up to 250 gallons per minute (gpm) or less, with occasional peak flows of approximately 500 gpm during significant precipitation events. Temporary dewatering will be conducted from wells located around the foundation excavation. Construction dewatering will include piping and discharging to a storm drain in the vicinity of the Site (BWSC catch basin at the southern end of Penniman Road) that discharges into the Charles River through outfall SDO037 (see Figure 3). Prior to discharge to remove suspended solids and dissolved and undissolved chemical constituents, collected water will be routed through a fractionation tank, bag filters and other necessary treatment components (i.e., granular activated carbon, GAC vessels) as shown on Figure 4. A Notice of Change (NOC) will be submitted to EPA if additional treatment components need to be mobilized at the Site.

DOCUMENTATION OF NATIONAL HISTORIC PRESERVATION ACT ELIGIBILITY REQUIREMENTS

Based on a review of the resources provided by the U.S. National Register of Historic Places and a review of the Massachusetts Cultural Resource Information System (MACRIS), no historic properties have been established to be present at the project Site, and discharges and discharge-related activities are not considered to have the potential to affect historic properties. The discharge is considered to meet Criterion A. Documentation is included in Appendix D.

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Reference: Remediation General Permit (RGP) 30 Penniman Road Allston, Massachusetts

DETERMINATION OF ENDANGERED SPECIES ACT ELIGIBILITY

According to the guidelines outlined in Appendix I of the 2017 NPDES RGP, a preliminary determination for the action area associated with this project was established using the U.S. Fish and Wildlife Service (FWS) Information for Planning and Consultation (IPaC) online system. A copy of the determination is attached in Appendix E. Based on the results of the determination, the project and action area are considered to meet FWS Criterion A as no listed species or critical habitat have been established to be present within the project action area.

SUPPLEMENTAL INFORMATION

An application for temporary Dewatering Discharge Permit Application is being submitted concurrently to the Boston Water and Sewer Commission. A copy this application is provided in Appendix F. Approval will be received prior to the start of discharge. The Department of Conservation and Recreation (DCR) has determined the stormwater system is not a DCR asset and does not require DCR permitting for its use (see Appendix F).

A Best Management Practices plan (BMP), which outlines the proposed discharge operations covered under the RGP, will be available at the site and the requirements are included in Appendix G.

OWNER AND OPERATOR INFORMATION

Owner and Operator:

30 Penn, LLC 675 VFW Pkwy #195 Chestnut Hill, Massachusetts 02467 Attn: Steve Ballas T: 617-888-3424 February 21, 2020 US Environmental Protection Agency Page 5 of 5

Reference: Remediation General Permit (RGP) 30 Penniman Road Allston, Massachusetts

CLOSING

Thank you very much for your consideration. Please feel free to contact us at your convenience.

Regards,

Stantec Consulting Services, Inc.

Juhan Jeannel
Richard Learned, LSP

Senior Environmental Project Manager

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Richard.Learned@stantec.com

Joseph Salvetti, LSP

Senior Associate Phone: 508.591.4327

Joseph.Salvetti@stantec.com

Attachment: FIGURES

TABLES

APPENDIX A – COPY OF NOTICE OF INTENT APPENDIX B – COPIES OF LABORATORY REPORTS

APPENDIX C – DILUTION FACTOR AND EFFLUENT LIMIT CALCULATIONS APPENDIX D – NATIONAL REGISTER OF HISTORIC PLACES DOCUMENTATION

APPENDIX E - ENDANGERED SPECIES ACT DOCUMENTATION

APPENDIX F - COPIES OF BWSC PERMIT APPLICATION AND DCR CORRESPONDENCE

APPENDIX G - BEST MANAGEMENT PRACTICES PLAN (BMPP)

c. 30 Penn, LLC MassDEP

Boston Water and Sewer Commission

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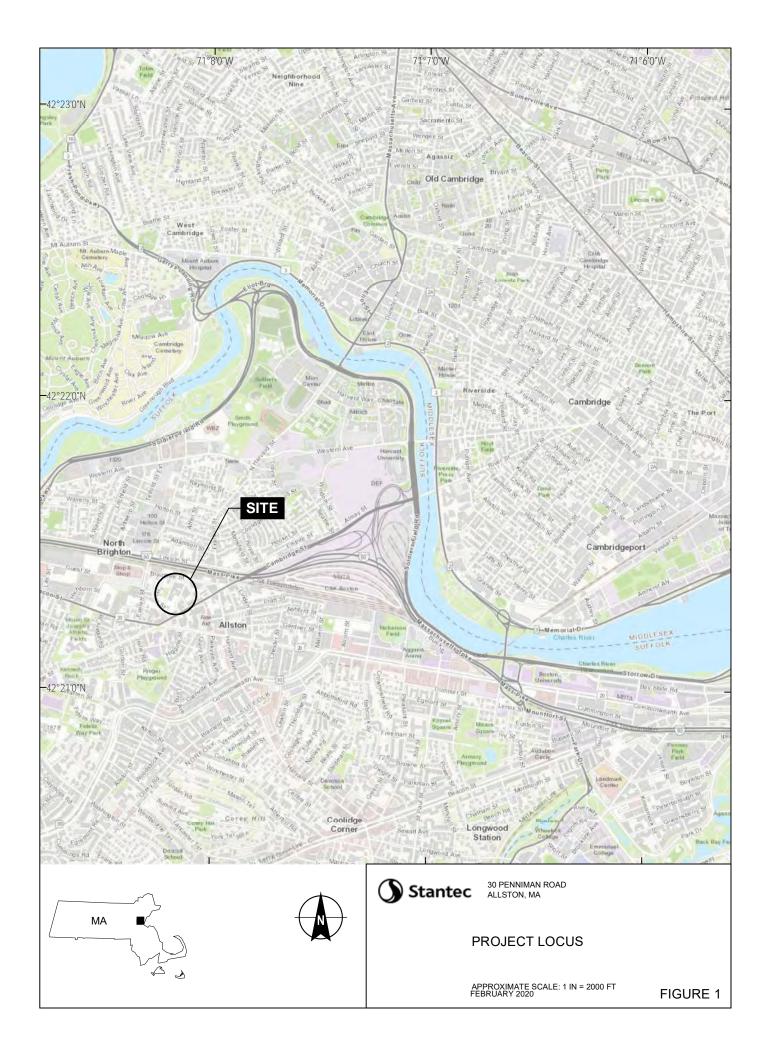


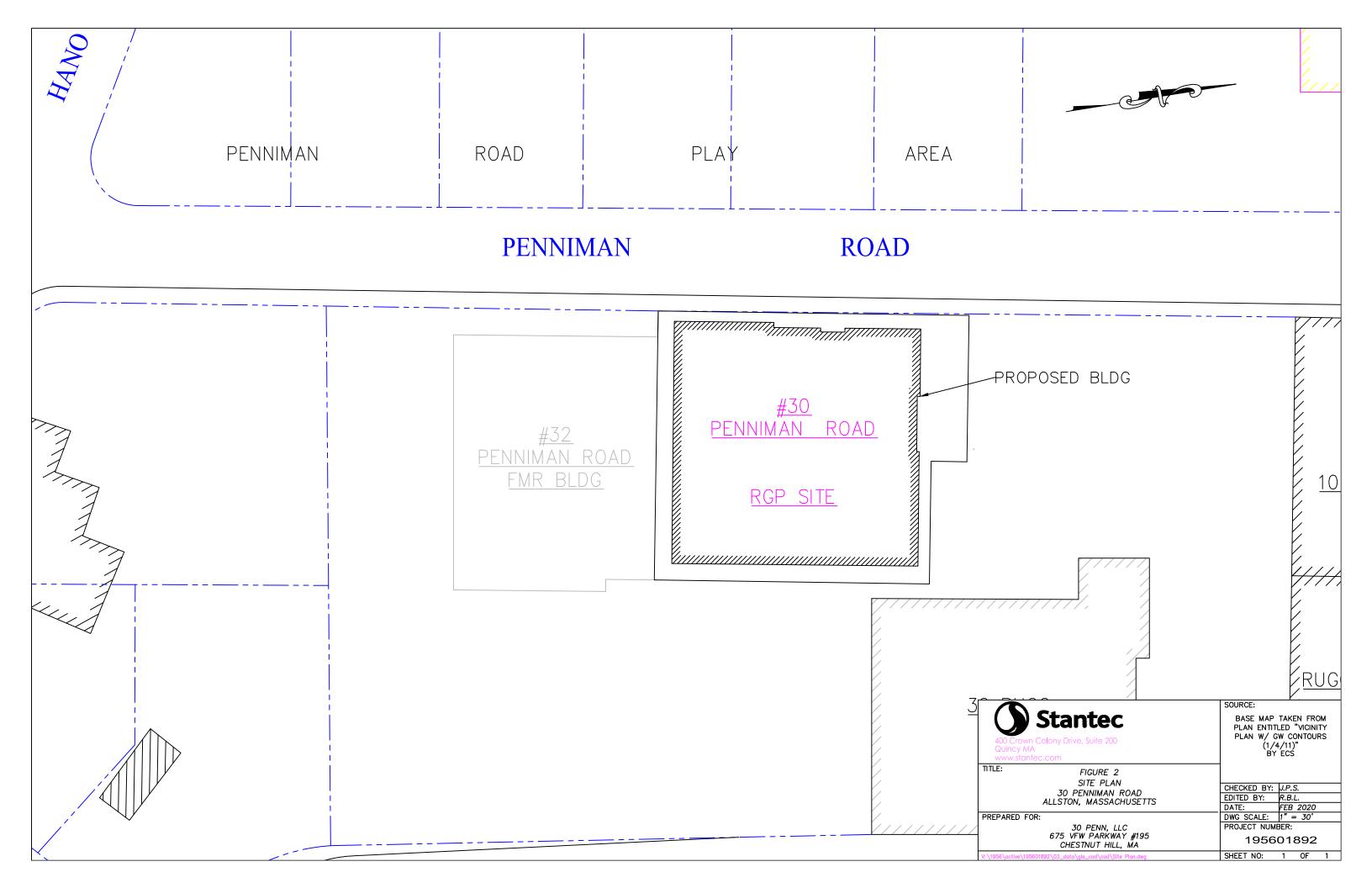
FIGURES

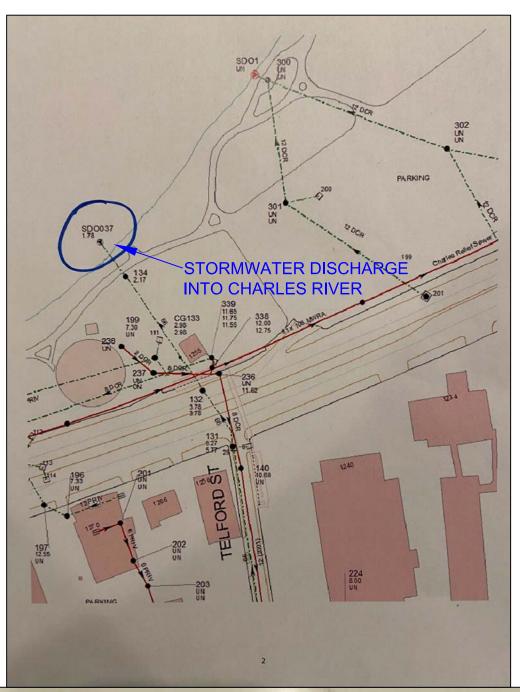
FIGURE 1 PROJECT LOCUS SITE PLAN

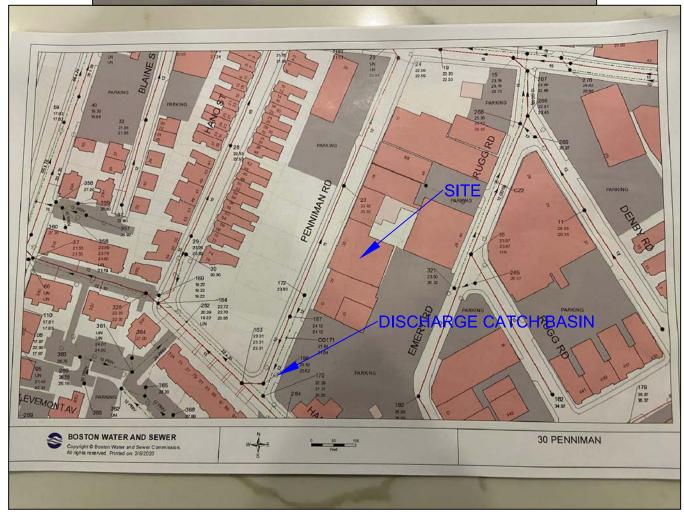
FIGURE 3 STORMWATER DISCHARGE PLAN

FIGURE 4 PROJECT TREATMENT SYSTEM SCHEMATIC











TITLE:

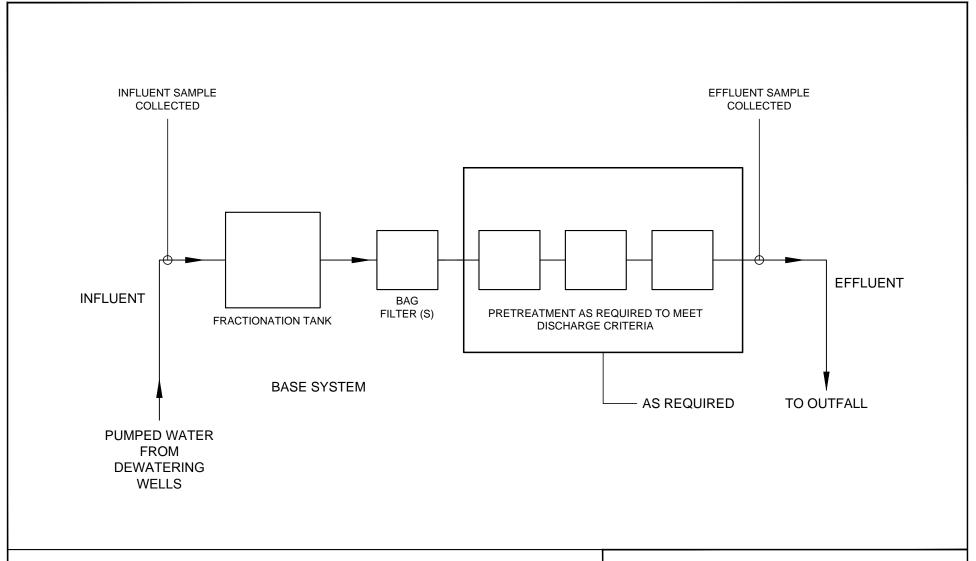
FIGURE 2 STORMWATER DISCHARGE PLAN 30 PENNIMAN ROAD ALLSTON, MASSACHUSETTS

PREPARED FOR:

30 PENN, LLC 675 VFW PARKWAY #195 CHESTNUT HILL, MA PLANS FROM BWSC FILES

CHECKED BY:	J.P.S.
EDITED BY:	R.B.L.
DATE:	FEB 2020
DWG SCALE:	not to scale
PROJECT NUMB	ER:

195601892 SHEET NO: 1 OF



LEGEND:



NOTE:

 DETAILS OF TREATMENT SYSTEM MAY VARY FROM SYSTEM INDICATED ABOVE. SPECIFIC MEANS AND METHODS OF TREATMENT TO BE SELECTED BY CONTRACTOR. WATER WILL BE TREATED TO MEET REQUIRED EFFLUENT STANDARDS.



30 PENNIMAN ROAD ALLSTON, MASSACHUSETTS

PROJECT TREATMENT SYSTEM SCHEMATIC

FEBRUARY 2020 FIGURE 4



TABLES

PARAMETERS IN GROUNDWATER FOR REMEDIATION GENERAL PERMIT (RGP) CRITERIA
PARAMETERS IN SURFACE WATER FOR REMEDIATION GENERAL PERMIT (RGP) CRITERIA

Phoenix Environmental Laboratories, Inc Table 1 - Parameters in Groundwater for Remediation General Permit (RGP) Criteria 587 East Middle Turnpike

587 East Middle Turnpike						
P.O. Box 370	Lab Sample Id			CF10994	CF10995	CF11044
Manchester, CT 06040	Collection Date			01/10/2020	01/10/2020	01/13/2020
(860) 645-1102	Client Id			C4	D1	TRIP BLANK
·	Matrix			GW	GW	WATER
Project: 30 PENNIMAN	CAS	Units	RCGW-2	Result	Result	Result
Miscellaneous/Inorganics						
1,2-Dibromoethane (EDB)	106-93-4	ug/L	2		< 0.02	
Ammonia as Nitrogen	7664-41-7	mg/L	10		0.7	
Chloride	16887-00-6	mg/L			291	
Chlorine Residual	7782-50-5	mg/L			< 0.02	
Ethanol	64-17-5	ug/L	10000		< 400	
O&G, Non-polar Material	PHNX - OIL-GREASE-NP	mg/L			< 1.4	
Total Cyanide	57-12-5	mg/L	0.03		< 0.010	
Total Suspended Solids	PHNX - TOTSUSPENDSOL	mg/L			56	
pH	field	pH Units		6.7	6.6	
Hardness	PHNX - HARDNESS	mg/L		567	470	
Metals Total	7440.00.0			.0.00=	.0.00=	
Antimony	7440-36-0	mg/L	8	< 0.005	< 0.005	
Arsenic	7440-38-2	mg/L	0.9	< 0.004	0.011	
Cadmium	7440-43-9	mg/L	0.004	< 0.001	< 0.001	
Chromium	7440-47-3	mg/L	0.3	0.009	0.007	
Chromium, Trivalent	16065-83-1	mg/L	0.6		0.007	
Chromium, Hexavalent	18540-29-9	mg/L	0.3		< 0.01	
Copper	7440-50-8	mg/L	100		0.005	
Iron	7439-89-6	mg/L			10	
Lead	7439-92-1	mg/L	0.01	0.006	0.009	
Mercury	7439-97-6	mg/L	0.02	< 0.0002	< 0.0002	
Nickel	7440-02-0	mg/L	0.2	0.008	0.011	
Selenium	7782-49-2	mg/L	0.1	< 0.010	< 0.010	
Silver	7440-22-4	mg/L	0.007	< 0.001	< 0.001	
Zinc	7440-66-6	mg/L	0.9	0.013	0.034	
Polychlorinated Biphenyls - SW8082A						
PCB-1016	12674-11-2	ug/L	5		< 0.048	
PCB-1221	11104-28-2	ug/L	5		< 0.048	
PCB-1232	11141-16-5	ug/L	5		< 0.048	
PCB-1242	53469-21-9	ug/L	5		< 0.048	
PCB-1248	12672-29-6	ug/L	5		< 0.048	
PCB-1254	11097-69-1	ug/L	5		< 0.048	
PCB-1260	11096-82-5	ug/L	5		< 0.048	
PCB-1262	37324-23-5	ug/L			< 0.048	
PCB-1268	11100-14-4	ug/L			< 0.048	
TOTAL PCBs					0	
MA Volotilo Detrolouro II udus soub sus (VDII)	VDU E /2004					
MA Volatile Petroleum Hydrocarbons (VPH) - MA C5-C8 Aliphatic Hydrocarbons *1,2	PHNX - C5-C8	/!	3000	< 100		
		ug/L				
C9-C12 Aliphatic Hydrocarbons *1,3	PHNX - C9-C12	ug/L	5000	< 100		
C9-C10 Aromatic Hydrocarbons *1	PHNX - C9-C10	ug/L	4000	< 100		
Benzene Ethyl Benzene	71-43-2	ug/L	1000	< 1.0		
Ethyl Benzene	100-41-4	ug/L	20000	< 1.0		·
m,p-Xylenes	179601-23-1	ug/L		< 2.0		
MTBE	1634-04-4	ug/L	50000	< 1.0		
Naphthalene	91-20-3	ug/L	700	< 5.0		
o-Xylene	95-47-6	ug/L		< 1.0		
Toluene	108-88-3	ug/L	50000	< 1.0		
TOTAL XYLENES			3000	0		
TOTAL BTEX				0		

Phoenix Environmental Laboratories, Inc Table 1 - Parameters in Groundwater for Remediation General Permit (RGP) Cri

P.O. Box 370	Lab Sample Id			CF10994	CF10995
Manchester, CT 06040 (860) 645-1102	Collection Date Client Id			01/10/2020 C4	01/10/20 D1
(800) 043-1102	Matrix			GW	GW
Project: 30 PENNIMAN	CAS	Units	RCGW-2	Result	Result
//A EPH Aliphatic/Aromatic Ranges - MAEPH 5/	2004				
9-C18 Aliphatic Hydrocarbons 1*	PHNX - C9-C18	ug/L	5000	< 190	
19-C36 Aliphatic Hydrocarbons 1*	PHNX - C19-C36	ug/L	50000	< 190	
11-C22 Aromatic Hydrocarbons 1,2* otal TPH 1,2*	PHNX - C11-C22 PHNX - EPH	ug/L ug/L	5000	< 190 < 190	
		- 01			
emivolatiles (SIM) - SW8270D (SIM)	83-32-9	ug/L	6000	0.64	< 0.48
Acenaphthylene	208-96-8	ug/L	40	0.13	< 0.10
Anthracene	120-12-7	ug/L	30	< 0.09	< 0.09
Benz(a)anthracene	56-55-3	ug/L	1000	< 0.09	< 0.10
Benzo(a)pyrene	50-32-8	ug/L	500	< 0.19	< 0.19
enzo(b)fluoranthene	205-99-2	ug/L	400	< 0.09	< 0.10
Benzo(ghi)perylene	191-24-2	ug/L	20	< 0.02	< 0.02
denzo(k)fluoranthene	207-08-9	ug/L	100 70	< 0.09	< 0.10
Chrysene Dibenz(a,h)anthracene	218-01-9 53-70-3	ug/L ug/L	40	< 0.05 < 0.02	< 0.05 < 0.02
luoranthene	206-44-0	ug/L ug/L	200	< 0.02	< 0.02
luorene	86-73-7	ug/L	40	1.1	< 0.10
lexachlorobenzene	118-74-1	ug/L	1		< 0.48
lexachlorobutadiene	87-68-3	ug/L	1		< 0.48
lexachlorocyclopentadiene	77-47-4	ug/L	5000		< 0.48
ndeno(1,2,3-cd)pyrene	193-39-5	ug/L	100	< 0.09	< 0.10
I-Nitrosodimethylamine	62-75-9	ug/L	5000	_	< 0.48
-Methylnaphthalene	91-57-6	ug/L	2000	< 0.47	< 0.48
laphthalene	91-20-3	ug/L	700	< 0.47	0.59
litrobenzene entachlorophenol	98-95-3 87-86-5	ug/L ug/L	50000 200		< 0.48
Phenanthrene	85-01-8	ug/L ug/L	10000	< 0.47	< 0.48
Pyrene	129-00-0	ug/L	20	0.1	< 0.45
yridine	110-86-1	ug/L	50000		< 1.9
TOTAL GROUP 1 PAI		<u> </u>		0	0
and alatha CMOSTOR					
emivolatiles - SW8270D ,2,4,5-Tetrachlorobenzene	95-94-3	ug/L	100000		< 3.3
,2,4-Trichlorobenzene	120-82-1	ug/L	200		< 4.8
,2-Dichlorobenzene	95-50-1	ug/L	2000		< 2.4
,2-Diphenylhydrazine	122-66-7	ug/L	5000		< 4.8
,3-Dichlorobenzene	541-73-1	ug/L	2000		< 2.4
,4-Dichlorobenzene	106-46-7	ug/L	200		< 2.4
,4,5-Trichlorophenol	95-95-4	ug/L	3000		< 0.95
,4,6-Trichlorophenol	88-06-2	ug/L	500		< 0.95
,4-Dichlorophenol	120-83-2	ug/L	2000		< 0.95
,4-Dimethylphenol ,4-Dinitrophenol	105-67-9 51-28-5	ug/L ug/L	40000 20000		< 0.95 < 0.95
,4-Dinitrophenoi	121-14-2	ug/L ug/L	20000		< 4.8
,6-Dinitrotoluene	606-20-2	ug/L	2000		< 4.8
-Chloronaphthalene	91-58-7	ug/L	100000		< 4.8
-Chlorophenol	95-57-8	ug/L	7000		< 0.95
-Methylphenol (o-cresol)	95-48-7	ug/L	50000		< 0.95
-Nitroaniline	88-74-4	ug/L			< 4.8
-Nitrophenol	88-75-5	ug/L	10000		< 0.95
&4-Methylphenol (m&p-cresol)	PHNX - M&P CRESOL	ug/L			< 9.5
,3'-Dichlorobenzidine	91-94-1	ug/L	2000		< 4.8
-Nitroaniline	99-09-2	ug/L	F000		< 4.8
,6-Dinitro-2-methylphenol	534-52-1	ug/L	5000		< 0.95
-Bromophenyl phenyl ether -Chloro-3-methylphenol	101-55-3 59-50-7	ug/L ug/L	10000		< 4.8 < 0.95
-Chloroaniline	106-47-8	ug/L ug/L	300		< 4.8
-Chlorophenyl phenyl ether	7005-72-3	ug/L ug/L	100000		< 0.95
-Nitroaniline	100-01-6	ug/L	100000		< 4.8
-Nitrophenol	100-02-7	ug/L	10000		< 0.95
cetophenone	98-86-2	ug/L	100000		< 4.8
niline	62-53-3	ug/L	100000		< 4.8
enzidine	92-87-5	ug/L	1000		< 4.8
enzoic acid	65-85-0	ug/L	100000		< 48
enzyl butyl phthalate	85-68-7	ug/L	10000		< 4.8
is(2-chloroethoxy)methane	111-91-1	ug/L	50000		< 4.8
is(2-chloroethyl)ether is(2-chloroisopropyl)ether	111-44-4 39638-32-9	ug/L ug/L	30 100		< 0.95 < 4.8
is(2-chioroisopropyi)ether is(2-ethylhexyl)phthalate	39638-32-9 117-81-7	ug/L ug/L	50000		< 4.8
arbazole	86-74-8	ug/L ug/L	50000		< 4.8
vi-n-butylphthalate	84-74-2	ug/L ug/L	5000		< 4.8
Di-n-octylphthalate	117-84-0	ug/L	100000		< 4.8
ibenzofuran	132-64-9	ug/L	10000		< 4.8
eiethyl phthalate	84-66-2	ug/L	9000		< 4.8
imethylphthalate	131-11-3	ug/L	50000		< 4.8
exachloroethane	67-72-1	ug/L	100		< 0.95
sophorone	78-59-1	ug/L	10000		< 4.8
l-Nitrosodi-n-propylamine	621-64-7	ug/L	5000		< 4.8
I-Nitrosodiphenylamine	86-30-6	ug/L	10000		< 4.8
entachloronitrobenzene	82-68-8	ug/L	10000		< 2.4
Phenol	108-95-2	ug/L	2000		< 0.95

Phoenix Environmental Laboratories, Inc. Table 1 - Parameters in Groundwater for Remediation General Permit (RGP) Crit 587 East Middle Turnpike

587 East Middle Turnpike	Lab Camanla Id			L 6510004	CE1000E
P.O. Box 370 Manchester, CT 06040	Lab Sample Id Collection Date			CF10994	CF10995 01/10/2020
(860) 645-1102	Client Id			C4	D1
	Matrix			GW	GW
Project: 30 PENNIMAN	CAS	Units	RCGW-2	Result	Result
Volatiles - SW8260C					
1,1,1,2-Tetrachloroethane	630-20-6	ug/L	10	< 1.0	< 1.0
1,1,1-Trichloroethane	71-55-6	ug/L	4000	< 1.0	< 1.0
1,1,2,2-Tetrachloroethane	79-34-5	ug/L	9	< 0.50	< 0.50
1,1,2-Trichloroethane	79-00-5	ug/L	900	< 1.0	< 1.0
1,1-Dichloroethane	75-34-3	ug/L	2000	< 1.0	< 1.0
1,1-Dichloroethene 1,1-Dichloropropene	75-35-4 563-58-6	ug/L ug/L	80	< 1.0 < 1.0	< 1.0 < 1.0
1,2,3-Trichlorobenzene	87-61-6	ug/L		< 1.0	< 1.0
1,2,3-Trichloropropane	96-18-4	ug/L	10000	< 1.0	< 1.0
1,2,4-Trichlorobenzene	120-82-1	ug/L	200	< 1.0	< 1.0
1,2,4-Trimethylbenzene	95-63-6	ug/L	100000	< 1.0	4.7
1,2-Dibromo-3-chloropropane	96-12-8	ug/L	1000	< 1.0	< 1.0
1,2-Dibromoethane 1,2-Dichlorobenzene	106-93-4 95-50-1	ug/L ug/L	2000	< 0.50 < 1.0	< 0.50 < 1.0
1,2-Dichloroethane	107-06-2	ug/L ug/L	5	< 0.60	< 0.60
1,2-Dichloropropane	78-87-5	ug/L	3	< 1.0	< 1.0
1,3,5-Trimethylbenzene	108-67-8	ug/L	1000	< 1.0	< 1.0
1,3-Dichlorobenzene	541-73-1	ug/L	2000	< 1.0	< 1.0
1,3-Dichloropropane	142-28-9	ug/L	50000	< 1.0	< 1.0
1,4-Dichlorobenzene	106-46-7	ug/L	200	< 1.0	< 1.0
2,2-Dichloropropane 2-Chlorotoluene	594-20-7 95-49-8	ug/L ug/L	10000	< 1.0 < 1.0	< 1.0 < 1.0
2-Hexanone	591-78-6	ug/L ug/L	10000	< 5.0	< 5.0
2-Isopropyltoluene	527-84-4	ug/L ug/L	10000	1.1	< 1.0
4-Chlorotoluene	106-43-4	ug/L		< 1.0	< 1.0
4-Methyl-2-pentanone	108-10-1	ug/L	50000	< 5.0	< 5.0
Acetone	67-64-1	ug/L	50000	< 25	< 25
Acrylonitrile	107-13-1	ug/L	10000	< 1.0	< 1.0
Benzene Bromobenzene	71-43-2 108-86-1	ug/L	1000 10000	< 0.70 < 1.0	< 0.70 < 1.0
Bromochloromethane	74-97-5	ug/L ug/L	10000	< 1.0	< 1.0
Bromodichloromethane	75-27-4	ug/L	6	< 0.50	< 0.50
Bromoform	75-25-2	ug/L	700	< 1.0	< 1.0
Bromomethane	74-83-9	ug/L	7	< 1.0	< 1.0
Carbon Disulfide	75-15-0	ug/L	10000	< 5.0	< 5.0
Carbon tetrachloride	56-23-5	ug/L	2	< 1.0	< 1.0
Chlorobenzene	108-90-7	ug/L	200	< 1.0	< 1.0
Chloroform Chloroform	75-00-3 67-66-3	ug/L	10000 50	< 1.0 < 1.0	< 1.0
Chloromethane	74-87-3	ug/L ug/L	10000	< 1.0	< 1.0 < 1.0
cis-1,2-Dichloroethene	156-59-2	ug/L	20	24	< 1.0
cis-1,3-Dichloropropene	10061-01-5	ug/L	5	< 0.40	< 0.40
Dibromochloromethane	124-48-1	ug/L	20	< 0.50	< 0.50
Dibromomethane	74-95-3	ug/L	50000	< 1.0	< 1.0
Dichlorodifluoromethane	75-71-8	ug/L	100000	< 1.0	< 1.0
Ethylbenzene Hexachlorobutadiene	100-41-4	ug/L	5000	< 1.0	1 .0.40
Isopropylbenzene	87-68-3 98-82-8	ug/L ug/L	100000	< 0.40	< 0.40
m&p-Xylene	179601-23-1	ug/L ug/L	100000	< 1.0	2.8
Methyl ethyl ketone	78-93-3	ug/L	5000	< 5.0	< 5.0
Methyl t-butyl ether (MTBE)	1634-04-4	ug/L	5000	< 1.0	< 1.0
Methylene chloride	75-09-2	ug/L	2000	< 1.0	< 1.0
n-Butylbenzene	104-51-8	ug/L		< 1.0	< 1.0
n-Propylbenzene	103-65-1	ug/L	10000	1.8	1.7
Naphthalene	91-20-3	ug/L	700	< 1.0	< 1.0
o-Xylene p-Isopropyltoluene	95-47-6 99-87-6	ug/L ug/L	6000 10000	< 1.0 < 1.0	< 1.0 < 1.0
sec-Butylbenzene	135-98-8	ug/L ug/L	10000	1.6	1.0
Styrene	100-42-5	ug/L	100	< 1.0	< 1.0
tert-Butylbenzene	98-06-6	ug/L	10000	< 1.0	< 1.0
Tetrachloroethene	127-18-4	ug/L	50	100	< 1.0
Tetrahydrofuran (THF)	109-99-9	ug/L	50000	< 2.5	< 2.5
Toluene Total Yulonos	108-88-3	ug/L	40000	< 1.0	< 1.0
Total Xylenes trans-1,2-Dichloroethene	1330-20-7 156-60-5	ug/L ug/L	500 80	< 1.0 5.4	2.8 < 1.0
trans-1,3-Dichloropropene	10061-02-6	ug/L ug/L	5	< 0.40	< 0.40
trans-1,4-dichloro-2-butene	110-57-6	ug/L	1000	< 5.0	< 5.0
Trichloroethene	79-01-6	ug/L	5	67	< 1.0
Trichlorofluoromethane	75-69-4	ug/L	100000	< 1.0	< 1.0
Trichlorotrifluoroethane	76-13-1	ug/L		< 1.0	< 1.0
Vinyl chloride	75-01-4	ug/L	2	10 212.9	< 1.0 15.9
TOTAL VOCs TOTAL BTEX				212.9	15.9 3.8
TOTAL BIEX					5.0
1,4-dioxane - SW8270DSIM					
1,4-dioxane	123-91-1	ug/l	6000		< 0.20
Oxygenates & Dioxane - SW8260C (OXY)					
1,4-Dioxane	123-91-1	ug/L	6000	< 50	< 50
Di-isopropyl ether Diethyl ether	108-20-3 60-29-7	ug/L ug/L	10000	< 1.0 < 1.0	< 1.0 < 1.0
Ethyl tert-butyl ether	637-92-3	ug/L ug/L	10000	< 1.0	< 1.0
tert-amyl methyl ether	994-05-8	ug/L		< 1.0	< 1.0
· · · ·		<u> </u>		<u> </u>	

Result Detected
Result Exceeds Criteria

Phoenix Environmental Laboratories, Inc. Table 2 - Parameters in Surface Water for Remediation General Permit (RGP) Criteria 587 East Middle Turnpike P.O. Box 370 Lab Sample Id CF32691 Manchester, CT 06040 **Collection Date** 2/14/2020 **CHARLES RIVER** (860) 645-1102 Client Id Matrix Surface Water Project Id: CHARLES RIVER CAS Units RL Result Miscellaneous/Inorganics 0.14 0.05 Ammonia as Nitrogen 7664-41-7 mg/L 1.00 рΗ PHNX - PH pH Units 7.31 0.5 Salinity 7647-14-5 0.5 ppt PHNX - HARDNESS 0.1 Hardness (CaCO3) 74.4 mg/L Metals, Total 7440-36-0 < 0.005 0.005 Antimony mg/L Arsenic 7440-38-2 < 0.004 0.004 mg/L < 0.001 Cadmium 7440-43-9 mg/L 0.001 Chromium 7440-47-3 mg/L < 0.001 0.001 0.01 Chromium, Trivalent 16065-83-1 mg/L < 0.01 0.01 Chromium, Hexavalent 18540-29-9 mg/L < 0.01 0.005 Copper 7440-50-8 mg/L < 0.005 Iron 7439-89-6 0.494 0.010 mg/L 0.002 Lead 7439-92-1 mg/L < 0.002 0.0002 7439-97-6 < 0.0002 Mercury mg/L

7440-02-0

7782-49-2

7440-22-4

7440-66-6

mg/L

mg/L

mg/L

mg/L

Notes

Nickel

Silver

Zinc

Selenium

RL - laboratory reporting limit

Result Detected

0.001

0.010

0.001

0.004

< 0.001

< 0.010

< 0.001

0.009



APPENDIX A

COPY OF NOTICE OF INTENT

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

A. General site information:

1. Name of site:	Site address: 30 PENNIMAN ROAD					
COMMERCIAL / RESIDENTIAL PROPERTY	Street:					
	City: ALLSTON		State: MA	Zip: 02134		
2. Site owner	Contact Person: STEPHEN BALLAS					
30 PENN, LLC	Telephone: 617.888.3424	Email: ST	EVE@BALI	LASGROUP.COM		
	Mailing address: 675 VFW PKWY #195					
	Street:					
Owner is (check one): ☐ Federal ☐ State/Tribal ■ Private ☐ Other; if so, specify:	City: CHESTNUT HILL	State: MA	Zip: 02467			
3. Site operator, if different than owner	Contact Person: STEPHEN BALLAS					
NA	Telephone: 617.888.3424	Email: ST	EVE@BALI	_ASGROUP.COM		
	Mailing address:					
	Street: SAME					
	City: SAME		State:	Zip:		
4. NPDES permit number assigned by EPA:	5. Other regulatory program(s) that apply to the site	(check all th	at apply):			
NA	■ MA Chapter 21e; list RTN(s): 3-36120	□ CERCI	LΑ			
		□ UIC Pro	ogram			
NPDES permit is (check all that apply: □ RGP □ DGP □ CGP	☐ NH Groundwater Management Permit or Groundwater Release Detection Permit:	\square POTW	Pretreatment	t		
☐ MSGP ☐ Individual NPDES permit ☐ Other; if so, specify:	Groundwater Release Detection Ferfilit.	□ CWA S	ection 404			

VIII? (check one):

■ Yes □ No

В.	Receiving	water	inforr	nation:

B. Receiving water information:			
1. Name of receiving water(s):	Waterbody identification of receiving water	(s): Classif	ication of receiving water(s):
CHARLES RIVER	MA72-36	CLASS	В
Receiving water is (check any that apply): □ Outstan	nding Resource Water □ Ocean Sanctuary □ territo	rial sea □ Wild and Scenic F	liver
2. Has the operator attached a location map in accord	dance with the instructions in B, above? (check one)	: ■ Yes □ No	
Are sensitive receptors present near the site? (check If yes, specify:	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. The Charles River at the outfall is li	s available for any of the indicated pollutants. For m	nore information, contact the	appropriate State as noted in Part
4. Indicate the seven day-ten-year low flow (7Q10) of Appendix V for sites located in Massachusetts and A		n the instructions in	15.70 MGD
5. Indicate the requested dilution factor for the calculaccordance with the instructions in Appendix V for s			22.81
6. Has the operator received confirmation from the a If yes, indicate date confirmation received: February 1. The street operator attached a summary of receiving (check one): ■ Yes □ No	12, 2020		
C. Source water information:			
1. Source water(s) is (check any that apply):			
■ Contaminated groundwater	☐ 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 than the receiving water; if	f
in accordance with the instruction in Appendix	RGP in accordance with the instruction in	so, indicate waterbody:	☐ Other; if so, specify:

Appendix VIII? (check one):

□ Yes □ No

2. Source water contaminants: PETROLEUM, METALS, VOCS, SVOCS	
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): \square Yes \blacksquare No If yes, indicate the contaminant(s) and	with the instructions in Appendix VIII? (check one): \square Yes \square No
the maximum concentration present in accordance with the instructions in	with the instructions in Appendix viii: (check one).
Appendix VIII.	
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): \square Existing discharge \blacksquare Nev	v discharge □ New source
Outfall(s):	Outfall location(s): (Latitude, Longitude)
SDO037	42.36527, -71.13859
Discharges enter the receiving water(s) via (check any that apply): □ Direct di	scharge to the receiving water ■ 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	var quatame
Has notification been provided to the owner of this system? (check one):	•
obtaining permission: BWSC permit being submitted concurrently. Appro	or discharges? (check one): ☐ Yes ■ No, if so, explain, with an estimated timeframe for
Has the operator attached a summary of any additional requirements the owner	·
Provide the expected start and end dates of discharge(s) (month/year): March	2020 to August 2020
Indicate if the discharge is expected to occur over a duration of: ■ less than 1	-
Has the operator attached a site plan in accordance with the instructions in D, a	DOVE! (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)
□ I – Petroleum-Related Site Remediation	 □ A. Inorganics □ B. Non-Halogenated Volatile Organi □ C. Halogenated Volatile Organic Cor □ D. Non-Halogenated Semi-Volatile Organi □ E. Halogenated Semi-Volatile Organi □ F. Fuels Parameters 	mpounds Organic Compounds
☐ II – Non-Petroleum-Related Site Remediation ■ III – Contaminated Site Dewatering		V, V, VI, VII or VIII: (check either G or H)
□ 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 □ 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 Kno	n Known		TF. 4		Influent		Effluent Limitations	
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		✓	1	E350.1	250	700	700	Report mg/L	
Chloride		✓	1	4500CLE+	15000	291000	291000	Report μg/l	
Total Residual Chlorine	✓		1	4500-CI-G	20	0	0	0.2 mg/L	
Total Suspended Solids		✓	1	2540D-11	10000	56000	56000	30 mg/L	
Antimony	✓		2	6010D	5	0	0	206 μg/L	
Arsenic		√	2	6010D	4	11	11	104 μg/L	
Cadmium		✓	2	6010D	1	0	0	10.2 μg/L	
Chromium III		✓	2	CALC	1	9	9	323 μg/L	
Chromium VI	✓		1	6010D	10	0	0	323 μg/L	
Copper		✓	1	6010D	5	5	5	242 μg/L	
Iron		✓	1	6010D	10	10000	10000	5,000 μg/L	
Lead		✓	2	6010D	2	9	9	160 μg/L	
Mercury	✓		2	7470A	0.2	0	0	0.739 μg/L	
Nickel		✓	2	6010D	1	11	11	1,450 μg/L	
Selenium	√		2	6010D	10	0	0	235.8 μg/L	
Silver	✓		2	6010D	1	0	0	35.1 μg/L	
Zinc		✓	2	6010D	4	34	34	420 μg/L	
Cyanide	✓		1	9010C/SW ₊	10	0	0	178 mg/L	
B. Non-Halogenated VOC									
Total BTEX	✓		2	8260C	1	3.8	3.8	100 μg/L	
Benzene	✓		2	8260C	0.7	0	0	5.0 μg/L	
1,4 Dioxane	✓		2	8260C	0.2	0	0	200 μg/L	
Acetone	✓		2	8260C	25	0	0	7.97 mg/L	
Phenol	✓		1	8260C	0.95	0	0	1,080 μg/L	

	Known	Known				Influent		Effluent Limitation	
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	✓		2	8260C	1	0	0	4.4 μg/L	
1,2 Dichlorobenzene	✓		2	8260C	1	0	0	600 μg/L	
1,3 Dichlorobenzene	✓		2	8260C	1	0	0	320 μg/L	
1,4 Dichlorobenzene	✓		2	8260C	1	0	0	5.0 μg/L	
Total dichlorobenzene	✓		2	8260C	1	0	0	763 μg/L in NH	
1,1 Dichloroethane	✓		2	8260C	1	0	0	70 μg/L	
1,2 Dichloroethane	✓		2	8260C	0.5	0	0	5.0 μg/L	
1,1 Dichloroethylene	✓		2	8260C	1	0	0	3.2 μg/L	
Ethylene Dibromide	✓		1	8260C	0.02	0	0	0.05 μg/L	
Methylene Chloride	✓		2	8260C	1	0	0	4.6 μg/L	
1,1,1 Trichloroethane	✓		2	8260C	1	0	0	200 μg/L	
1,1,2 Trichloroethane	✓		2	8260C	1	0	0	5.0 μg/L	
Trichloroethylene		✓	2	8260C	10	67	67	5.0 μg/L	
Tetrachloroethylene		✓	2	8260C	10	100	100	5.0 μg/L	
cis-1,2 Dichloroethylene		✓	2	8260C	10	24	24	70 μg/L	
Vinyl Chloride		✓	2	8260C	1	10	10	2.0 μg/L	
D. Non-Halogenated SVOC	Cs.								
Total Phthalates	✓		1	8270D	4.8	0	0	190 μg/L	
Diethylhexyl phthalate	✓		1	8270D	4.8	0	0	101 μg/L	
Total Group I PAHs	✓		1	8270D	0.48	0	0	1.0 μg/L	
Benzo(a)anthracene	✓		2	8270D	0.09	0	0		
Benzo(a)pyrene	✓		2	8270D	0.19	0	0		
Benzo(b)fluoranthene	✓		2	8270D	0.09	0	0		
Benzo(k)fluoranthene	√		2	8270D	0.09	0	0	As Total PAHs	
Chrysene	✓		2	8270D	0.05	0	0	1	
Dibenzo(a,h)anthracene	✓		2	8270D	0.02	0	0	1	
Indeno(1,2,3-cd)pyrene	✓		2	8270D	0.09	0	0	†	

	Known	Known		Test method (#)		In	fluent	Effluent Limitations		
Parameter	or believed absent	or believed present	# of samples		Detection limit (µg/l)	Daily maximum (µg/l)	Daily average (µg/l)	TBEL	WQBEL	
Total Group II PAHs	✓		2	8270D	VARIES	0	0	100 μg/L		
Naphthalene	✓		2	8270D	0.47	0	0	20 μg/L		
E. Halogenated SVOCs										
Total PCBs	✓		1	8082A	0.048	0	0	0.000064 μg/L		
Pentachlorophenol	✓		1	8270D	0.048	0	0	1.0 μg/L		
F. Fuels Parameters										
Total Petroleum Hydrocarbons	✓		2	E1664A	0.0014	0	0	5.0 mg/L		
Ethanol	✓		1	8260C	400	0	0	Report mg/L		
Methyl-tert-Butyl Ether	✓		2	8260C	1	0	0	70 μg/L		
tert-Butyl Alcohol	✓		1	8260C	1	0	0	120 μg/L in MA 40 μg/L in NH		
tert-Amyl Methyl Ether	✓		1	8260C	1	0	0	90 μg/L in MA 140 μg/L in NH		
Other (i.e., pH, temperatur	e, hardness,	salinity, LC	C ₅₀ , addition	nal pollutar	nts present);	if so, specify:	6.6			
Hardness (mg/L)		✓	1	Lab		567	567			
Temperature (deg C)		✓	1	field		7.7	7.7			
Salinity (ppt)	✓		1	field		0.5	0.5			
	1									

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. Groundwater pumped from a ring of extraction wells will be pumped to dewater the excavation for construction of building footings and elevator shaft. The pumped water collected in a Frac tank, pumped through sediment filters then through granular activated carbon (GAC) vessels prior to discharge into the Boston Waster and Sewer Communication at a catch basin located at the southern end of Penniman Road in Allston, MA. This system discharges to the Charles River at outfall SDO037. 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: GAC vessels 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: Treatment system Is use of a flow meter feasible? (check one): ■ Yes □ No, if so, provide justification:	500
Provide the proposed maximum effluent flow in gpm.	500
Provide the average effluent flow in gpm.	250
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

Seavengers 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): Yes 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 operat	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)
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)). 8. 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): □ Yes □ 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	□ Algaecides/biocides □ Antifoams □ Coagulants □ Corrosion/scale inhibitors □ Disinfectants □ Flocculants □ Neutralizing agents □ Oxidants □ Oxygen □
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): G. Part 423.15(j)(1) are non-detect in discharges with the addition of the proposed chemical/additive? (check one): 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 FWS Criterion C: Using the best scientific and commercial data available, the effect of the discharges and related activities on listed species and critical	scavengers □ pH conditioners □ Bioremedial agents, including microbes □ Chlorine or chemicals containing chlorine □ Other; if so, specify:
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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	
☐ FWS Criterion C: Using the best scientific and commercial data available, the effect of the discharges and related activities on listed species and critical	(informal consultation). Has the operator completed consultation with FWS? (check one): ☐ Yes ☐ No; if no, is consultation underway? (check one): ☐
	Yes □ No
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) \Box the operator \Box EPA \Box Other; if so, specify:	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

□ 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): ■ Yes □ 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): Yes 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

Print Name and Title: Stephen Ballas, Manager

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 assure that qualified personnel properly gathered and evaluated the information submitted. Based on my inquiry of the person or persons who manage the system, or those persons directly responsible for gathering the information, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I have no personal knowledge that the information submitted is other 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.										
A BMP PLAN WILL BE IMPLEMENTED PRIOR TO START OF DISC BMPP certification statement:	CHARGE									
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										
$permit(s). \ Additional \ discharge \ permit \ is \ (check \ one): \ \Box \ RGP \ \Box \ DGP \ \Box \ CGP \ \Box \ MSGP \ \ \Box \ Individual \ NPDES \ permit$	Check one: Yes □ No □ NA ■									
☐ Other; if so, specify:										
Signature: Stephen Ballas, Manager Da	te: 02/21/2020									



APPENDIX B
COPIES OF LABORATORY REPORTS



Tuesday, January 21, 2020

Attn: Mr Richard Learned

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Project ID: 30 PENNIMAN SDG ID: GCF10994 Sample ID#s: CF10994

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,

Phyllis/Shiller

Laboratory Director

NELAC - #NY11301

CT Lab Registration #PH-0618

MA Lab Registration #M-CT007

ME Lab Registration #CT-007

NH Lab Registration #213693-A,B

NJ Lab Registration #CT-003 NY Lab Registration #11301

PA Lab Registration #68-03530

RI Lab Registration #63

UT Lab Registration #CT00007

VT Lab Registration #VT11301



Environmental Laboratories, Inc.

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



SDG Comments

January 21, 2020

SDG I.D.: GCF10994

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.



Environmental Laboratories, Inc.

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

Sample Id Cross Reference

January 21, 2020

SDG I.D.: GCF10994

Project ID: 30 PENNIMAN

Client Id	Lab Id	Matrix
C4	CF10994	GROUND WATER



Environmental Laboratories, Inc.

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

Analysis Report

January 21, 2020

FOR: Attn: Mr Richard Learned

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample InformationCustody InformationDateTimeMatrix:GROUND WATERCollected by:01/10/2013:00Location Code:STANTECMAReceived by:CP01/13/2013:30

Rush Request: Standard Analyzed by: see "By" below

Laboratory Data

SDG ID: GCF10994
Phoenix ID: CF10994

Project ID: 30 PENNIMAN

195601892

Client ID: C4

P.O.#:

	5	RL/		5	D . (T)	_	5 /
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	01/15/20	EK	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	01/15/20	EK	SW6010D
Barium	0.094	0.002	mg/L	1	01/15/20	EK	SW6010D
Beryllium	< 0.001	0.001	mg/L	1	01/15/20	EK	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	01/15/20	EK	SW6010D
Chromium	0.009	0.001	mg/L	1	01/15/20	EK	SW6010D
Mercury	< 0.0002	0.0002	mg/L	1	01/15/20	RS	SW7470A
Nickel	0.008	0.001	mg/L	1	01/15/20	EK	SW6010D
Lead	0.006	0.002	mg/L	1	01/15/20	EK	SW6010D
Antimony	< 0.005	0.005	mg/L	1	01/15/20	EK	SW6010D
Selenium	< 0.010	0.010	mg/L	1	01/15/20	EK	SW6010D
Thallium	< 0.0005	0.0005	mg/L	5	01/14/20	CPP	SW6020B
Vanadium	0.019	0.002	mg/L	1	01/15/20	EK	SW6010D
Zinc	0.013	0.004	mg/L	1	01/15/20	EK	SW6010D
Mercury Digestion	Completed				01/15/20	Q/Q	SW7470A
EPH Extraction	Completed				01/17/20	JS/JS	SW3510C
MA Petroleum Hydrocarbon (EPH)	Completed				01/13/20		MADEP EPH-04
Semi-Volatile Extraction	Completed				01/13/20	P/AK	SW3520C
Total Metals Digestion	Completed				01/14/20	AG	
Total Metals Digestion MS	Completed				01/13/20	AG	
MA Petroleum Hydrocarbon (VPH)	Completed				01/13/20	RM	MADEP VPH04
<u>Volatiles</u>							
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,1,1-Trichloroethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	01/14/20	MH	SW8260C
1,1,2-Trichloroethane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,1-Dichloroethane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C

Client ID: C4

Client ID. C4		D1 /					
Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
1,1-Dichloroethene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,1-Dichloropropene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2,3-Trichloropropane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2-Dibromoethane	ND	0.50	ug/L	1	01/14/20	MH	SW8260C
1,2-Dichlorobenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
1,2-Dichloroethane	ND	0.60	ug/L	1	01/14/20	МН	SW8260C
1,2-Dichloropropane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,3-Dichlorobenzene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,3-Dichloropropane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
1,4-Dichlorobenzene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
2,2-Dichloropropane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
2-Chlorotoluene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
2-Hexanone	ND	5.0	ug/L	1	01/14/20	МН	SW8260C
2-Isopropyltoluene	1.1	1.0	ug/L	1	01/14/20	МН	SW8260C
4-Chlorotoluene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
4-Methyl-2-pentanone	ND	5.0	ug/L	1	01/14/20	МН	SW8260C
Acetone	ND	25	ug/L	1	01/14/20	МН	SW8260C
Acrylonitrile	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
Benzene	ND	0.70	ug/L	1	01/14/20	MH	SW8260C
Bromobenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Bromochloromethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Bromodichloromethane	ND	0.50	ug/L	1	01/14/20	MH	SW8260C
Bromoform	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
Bromomethane	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
Carbon Disulfide	ND	5.0	ug/L	1	01/14/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
cis-1,2-Dichloroethene	24	1.0	ug/L	10	01/14/20	MH	SW8260C
,	ND	0.40	ug/L	10	01/13/20	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40		1	01/14/20	МН	SW8260C SW8260C
Dibromochloromethane			ug/L	1			SW8260C SW8260C
Dibromomethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C SW8260C
Dichlorodifluoromethane	ND	1.0	ug/L	1	01/14/20	MH	
Ethylbenzene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	01/14/20	MH	SW8260C
Isopropylbenzene	2.0	1.0	ug/L	1	01/14/20	MH	SW8260C
m&p-Xylene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	01/14/20	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Methylene chloride	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Naphthalene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C

Project ID: 30 PENNIMAN Phoenix I.D.: CF10994

Client ID: C4

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
n-Propylbenzene	1.8	1.0	ug/L	1	01/14/20	МН	SW8260C
o-Xylene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
p-lsopropyltoluene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
sec-Butylbenzene	1.6	1.0	ug/L	1	01/14/20	МН	SW8260C
Styrene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	01/14/20	МН	SW8260C
Tetrachloroethene	100	10	ug/L	10	01/15/20	МН	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	01/14/20	MH	SW8260C
Toluene	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Total Xylenes	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
trans-1,2-Dichloroethene	5.4	1.0	ug/L	1	01/14/20	MH	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	01/14/20	MH	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	01/14/20	MH	SW8260C
Trichloroethene	67	10	ug/L	10	01/15/20	MH	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Trichlorotrifluoroethane	ND	1.0	ug/L	1	01/14/20	MH	SW8260C
Vinyl chloride	10	1.0	ug/L	1	01/14/20	MH	SW8260C
QA/QC Surrogates							
% 1,2-dichlorobenzene-d4	93		%	1	01/14/20	МН	70 - 130 %
% Bromofluorobenzene	96		%	1	01/14/20	MH	70 - 130 %
% Dibromofluoromethane	96		%	1	01/14/20	МН	70 - 130 %
% Toluene-d8	114		%	1	01/14/20	МН	70 - 130 %
% 1,2-dichlorobenzene-d4 (10x)	95		%	10	01/15/20	МН	70 - 130 %
% Bromofluorobenzene (10x)	95		%	10	01/15/20	МН	70 - 130 %
% Dibromofluoromethane (10x)	102		%	10	01/15/20	МН	70 - 130 %
% Toluene-d8 (10x)	98		%	10	01/15/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	01/14/20	MH	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	01/14/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	01/14/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	01/14/20	МН	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	01/14/20	MH	SW8260C (OXY)
Semivolatiles by SIM, PA	<u> </u>						
2-Methylnaphthalene	ND	0.47	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Acenaphthene	0.64	0.47	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Acenaphthylene	0.13	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Anthracene	ND	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Benz(a)anthracene	ND	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Benzo(a)pyrene	ND	0.19	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Benzo(b)fluoranthene	ND	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Benzo(ghi)perylene	ND	0.02	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Benzo(k)fluoranthene	ND	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Chrysene	ND	0.05	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Dibenz(a,h)anthracene	ND	0.02	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Fluoranthene	ND	0.47	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Fluorene	1.1	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Indeno(1,2,3-cd)pyrene	ND	0.09	ug/L	1	01/14/20	KCA	SW8270D (SIM)
Naphthalene	ND	0.47	ug/L	1	01/14/20	KCA	SW8270D (SIM)

Client ID: C4

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference	
Phenanthrene	ND	0.47	ug/L	1	01/14/20	KCA	SW8270D (SIM)	
Pyrene	0.10	0.07	ug/L	1	01/14/20	KCA	SW8270D (SIM)	
QA/QC Surrogates								
% 2-Fluorobiphenyl	47		%	1	01/14/20	KCA	40 - 140 %	
% Nitrobenzene-d5	42		%	1	01/14/20	KCA	40 - 140 %	
% Terphenyl-d14	36		%	1	01/14/20	KCA	40 - 140 %	3
MA EPH Aliphatic/Aroma	tic Rang	<u>jes</u>						
C11-C22 Aromatic Hydrocarbons 1,2*	ND	190	ug/L	1	01/17/20	AW	MAEPH 5/2004	
C11-C22 Aromatic Hydrocarbons Unadj	ND	190	ug/L	1	01/17/20	AW	MAEPH 5/2004	
C19-C36 Aliphatic Hydrocarbons 1*	ND	190	ug/L	1	01/18/20	AW	MAEPH 5/2004	
C9-C18 Aliphatic Hydrocarbons 1*	ND	190	ug/L	1	01/18/20	AW	MAEPH 5/2004	
Total TPH 1,2*	ND	190	ug/L	1	01/17/20	AW	MAEPH 5/2004	
QA/QC Surrogates								
% 1-chlorooctadecane (aliphatic)	22		%	1	01/18/20	AW	40 - 140 %	3
% 2-Bromonaphthalene (Fractionation)	71		%	1	01/17/20	AW	40 - 140 %	
% 2-Fluorobiphenyl (Fractionation)	74		%	1	01/17/20	AW	40 - 140 %	
% o-terphenyl (aromatic)	40		%	1	01/17/20	AW	40 - 140 %	
MA Volatile Petroleum H	ydrocark	ons (VPF	<u>l)</u>					
Unadjusted C5-C8 Aliphatics (*1)	ND	100	ug/L	1	01/13/20	RM	MA VPH 5/2004	
Unadjusted C9-C12 Aliphatics (*1)	ND	100	ug/L	1	01/13/20	RM	MA VPH 5/2004	
C5-C8 Aliphatic Hydrocarbons *1,2	ND	100	ug/L	1	01/13/20	RM	MA VPH 5/2004	
C9-C12 Aliphatic Hydrocarbons *1,3	ND	100	ug/L	1	01/13/20	RM	MA VPH 5/2004	
C9-C10 Aromatic Hydrocarbons *1	ND	100	ug/L	1	01/13/20	RM	MA VPH 5/2004	
Benzene	ND	1.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
Ethyl Benzene	ND	1.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
MTBE	ND	1.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
Naphthalene	ND	5.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
Toluene	ND	1.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
m,p-Xylenes	ND	2.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
o-Xylene	ND	1.0	ug/L	1	01/13/20	RM	MA VPH 5/2004	
QA/QC Surrogates								
% 2,5-Dibromotoluene (FID)	105		%	1	01/13/20	RM	70 - 130 %	
% 2,5-Dibromotoluene (PID)	101		%	1	01/13/20	RM	70 - 130 %	

Phoenix I.D.: CF10994

Project ID: 30 PENNIMAN Phoenix I.D.: CF10994

Client ID: C4

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

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:

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.

8260 Analysis:

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

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

EPH Comment

Poor Surrogate recovery. Insufficient sample for rextraction.

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

January 21, 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

QA/QC Data

January 21, 2020		QA/QC Data					SDG I.D.: GCF10994						
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 514110 (mg/L), Q	C Sam	ple No: (CF11603	(CF1099	94)								
Mercury - Water			<0.0002		NC	110			103			75 - 125	30
Comment:													
Additional Mercury criteria: LCS ac	ceptano	e range f	or waters	is 80-1209	% and fo	or soils is	s 75-1259	%					
QA/QC Batch 514163 (mg/L), Q	C Sam _l	ple No: (CF10467	(CF1099	94)								
ICP Metals - Aqueous													
Antimony	BRL	0.005	< 0.005	< 0.005	NC	102	104	1.9	114			75 - 125	20
Arsenic	BRL	0.004	< 0.004	< 0.004	NC	96.4	99.1	2.8	109			75 - 125	20
Barium	BRL	0.002	0.127	0.126	0.80	98.9	103	4.1	96.4			75 - 125	20
Beryllium	BRL	0.001	< 0.001	< 0.001	NC	98.6	101	2.4	96.8			75 - 125	20
Cadmium	BRL	0.001	0.004	0.004	NC	97.7	101	3.3	90.6			75 - 125	20
Chromium	BRL	0.001	< 0.001	< 0.001	NC	97.9	101	3.1	94.5			75 - 125	20
Lead	BRL	0.002	0.003	0.003	NC	94.1	96.9	2.9	88.4			75 - 125	20
Nickel	BRL	0.001	0.100	0.100	0	100	103	3.0	91.4			75 - 125	20
Selenium	BRL	0.010	0.012	0.012	NC	94.0	96.1	2.2	101			75 - 125	20
Silver	BRL	0.001	< 0.001	0.002	NC	95.0	97.8	2.9	112			75 - 125	20
Vanadium	BRL	0.002	0.091	0.092	1.10	96.6	99.2	2.7	94.9			75 - 125	20
Zinc	BRL	0.004	0.292	0.292	0	96.9	100	3.1	104			75 - 125	20
QA/QC Batch 513984 (mg/L), Q	C Sam _l	ple No: (CF10473	5X (CF1	0994)								
ICP MS Metals - Aqueous	<u>.</u>												
Thallium	BRL	0.0005	<0.0005	<0.0005	NC	105	107	1.9	103			75 - 125	20



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

QA/QC Report

QA/QC Data

January 21, 2020		<u>QA/C</u>	OC Data				SDG I	.D.: 0	GCF109	994	
Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
QA/QC Batch 514708 (ug/L), QC	Samp	le No: CF10994 (CF10994	.)								
MAEPH - Ground Water		•	•								
C9-C18 Aliphatic Hydrocarbons 1*	ND	100	61	53	14.0				40 - 140	25	
C19-C36 Aliphatic Hydrocarbons 1*	ND	100	70	63	10.5				40 - 140	25	
C11-C22 Aromatic Hydrocarbons 1	ND	100	64	60	6.5				40 - 140	25	
C11-C22 Aromatic Hydrocarbons U	ND	100	0.1	00	0.0				40 - 140	25	
Total TPH 1,2*	ND	100	66	58	12.9				40 - 140	25	
C9 - Nonane	ND	10	43	36	17.7				40 - 140	25	1
C-10 Decane	ND	10	54	46	16.0				40 - 140	25	•
C12 - Dodecane	ND	10	54	48	11.8				40 - 140	25	
C14 - Tetradecane	ND	10	61	55	10.3				40 - 140	25	
C16 - Hexadecane	ND	10	74	64	14.5				40 - 140	25	
C18 - Octadecane	ND	10	79	69	13.5				40 - 140	25	
C19 - Nonadecane	ND	10	77	67	13.9				40 - 140	25	
C20 - Eicosane	ND	10	77	67	13.9				40 - 140	25	
C22 - Docosane	ND	10	75	65	14.3				40 - 140	25	
C24 - Tetracosane	ND	10	77	67	13.9				40 - 140	25	
C26 - Hexacosane	ND	10	76	66	14.1				40 - 140	25	
C28 - Octacosane	ND	10	75	66	12.8				40 - 140	25	
C30 - Tricotane	ND	10	82	79	3.7				40 - 140	25	
C36 - Hexatriacontane	ND	10	23	25	8.3				40 - 140	25	1
% 1-chlorooctadecane (aliphatic)	45	%	57	55	3.6				40 - 140	25	
% o-terphenyl (aromatic)	60	%	59	58	1.7				40 - 140	25	
% 2-Fluorobiphenyl (Fractionation)	71	%	73	78	6.6				40 - 140	25	
% 2-Bromonaphthalene (Fractionati	68	%	73	81	10.4				40 - 140	25	
% 2-Methylnaphthalene BT		%	0	0	NC				0 - 5		
% Naphthalene BT		%	0	0	NC				0 - 5		
Comment:											
Additional EPH fractionation criteria	a: Break	through criteria (BT) is 0 to 5%	6								
QA/QC Batch 513953 (ug/L), QC		=									
Semivolatiles by SIM, PAF			,								
2-Methylnaphthalene	ND	0.50	59	54	8.8				40 - 140	20	
Acenaphthene	ND	0.50	65	63	3.1				40 - 140	20	
Acenaphthylene	ND	0.10	66	62	6.3				40 - 140	20	
Anthracene	ND	0.10	87	86	1.2				40 - 140	20	
Benz(a)anthracene	ND	0.02	92	88	4.4				40 - 140	20	
Benzo(a)pyrene	ND	0.02	85	78	8.6				40 - 140	20	
Benzo(b)fluoranthene	ND	0.02	81	74	9.0				40 - 140	20	
Benzo(ghi)perylene	ND	0.02	77	71	8.1				40 - 140	20	
Benzo(k)fluoranthene	ND	0.02	85	78	8.6				40 - 140	20	
Chrysene	ND	0.02	82	79	3.7				40 - 140	20	
Dibenz(a,h)anthracene	ND	0.02	81	75	7.7				40 - 140	20	
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SDG I.D.: GCF10994

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
Fluoranthene	ND	0.50	80	78	2.5				40 - 140	20
Fluorene	ND	0.10	71	70	1.4				40 - 140	20
Indeno(1,2,3-cd)pyrene	ND	0.02	75	70	6.9				40 - 140	20
Naphthalene	ND	0.50	48	45	6.5				40 - 140	20
Phenanthrene	ND	0.50	82	81	1.2				40 - 140	20
Pyrene	ND	0.07	83	81	2.4				40 - 140	20
% 2-Fluorobiphenyl	57	%	54	51	5.7				40 - 140	20
% Nitrobenzene-d5	48	%	47	43	8.9				40 - 140	20
% Terphenyl-d14	67	%	69	65	6.0				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 514113 (ug/L), QC Sample No: CF10770 (CF10994)

QA/QC Batch 514113 (ug/L), C	C Samp	le No: CF10770 (CF10994)									
Volatiles - Ground Water											
1,1,1,2-Tetrachloroethane	ND	1.0	101	107	5.8	106	106	0.0	70 - 130	30	
1,1,1-Trichloroethane	ND	1.0	96	105	9.0	105	105	0.0	70 - 130	30	
1,1,2,2-Tetrachloroethane	ND	0.50	100	108	7.7	105	105	0.0	70 - 130	30	
1,1,2-Trichloroethane	ND	1.0	97	104	7.0	94	96	2.1	70 - 130	30	
1,1-Dichloroethane	ND	1.0	97	104	7.0	102	102	0.0	70 - 130	30	
1,1-Dichloroethene	ND	1.0	101	111	9.4	112	112	0.0	70 - 130	30	
1,1-Dichloropropene	ND	1.0	95	102	7.1	112	108	3.6	70 - 130	30	
1,2,3-Trichlorobenzene	ND	1.0	78	101	25.7	88	108	20.4	70 - 130	30	
1,2,3-Trichloropropane	ND	1.0	91	95	4.3	90	95	5.4	70 - 130	30	
1,2,4-Trichlorobenzene	ND	1.0	86	104	18.9	97	109	11.7	70 - 130	30	
1,2,4-Trimethylbenzene	ND	1.0	96	104	8.0	103	100	3.0	70 - 130	30	
1,2-Dibromo-3-chloropropane	ND	1.0	93	110	16.7	101	114	12.1	70 - 130	30	
1,2-Dibromoethane	ND	1.0	99	103	4.0	100	102	2.0	70 - 130	30	
1,2-Dichlorobenzene	ND	1.0	96	103	7.0	100	101	1.0	70 - 130	30	
1,2-Dichloroethane	ND	1.0	98	104	5.9	101	102	1.0	70 - 130	30	
1,2-Dichloropropane	ND	1.0	96	103	7.0	107	106	0.9	70 - 130	30	
1,3,5-Trimethylbenzene	ND	1.0	96	104	8.0	104	102	1.9	70 - 130	30	
1,3-Dichlorobenzene	ND	1.0	95	101	6.1	100	100	0.0	70 - 130	30	
1,3-Dichloropropane	ND	1.0	97	103	6.0	96	97	1.0	70 - 130	30	
1,4-Dichlorobenzene	ND	1.0	93	103	10.2	100	101	1.0	70 - 130	30	
1,4-dioxane	ND	100	106	114	7.3	105	101	3.9	40 - 160	30	
2,2-Dichloropropane	ND	1.0	90	100	10.5	90	90	0.0	70 - 130	30	
2-Chlorotoluene	ND	1.0	98	105	6.9	104	104	0.0	70 - 130	30	
2-Hexanone	ND	5.0	85	95	11.1	89	94	5.5	40 - 160	30	
2-Isopropyltoluene	ND	1.0	99	108	8.7	108	107	0.9	70 - 130	30	
4-Chlorotoluene	ND	1.0	93	101	8.2	100	99	1.0	70 - 130	30	
4-Methyl-2-pentanone	ND	5.0	89	97	8.6	46	47	2.2	40 - 160	30	
Acetone	ND	5.0	105	96	9.0	93	107	14.0	40 - 160	30	
Acrylonitrile	ND	5.0	99	106	6.8	99	107	7.8	70 - 130	30	
Benzene	ND	0.70	95	102	7.1	108	106	1.9	70 - 130	30	
Bromobenzene	ND	1.0	99	104	4.9	102	103	1.0	70 - 130	30	
Bromochloromethane	ND	1.0	99	108	8.7	103	102	1.0	70 - 130	30	
Bromodichloromethane	ND	0.50	98	106	7.8	106	106	0.0	70 - 130	30	
Bromoform	ND	1.0	97	103	6.0	98	100	2.0	70 - 130	30	
Bromomethane	ND	1.0	100	113	12.2	80	87	8.4	40 - 160	30	
Carbon Disulfide	ND	1.0	92	101	9.3	99	101	2.0	70 - 130	30	
Carbon tetrachloride	ND	1.0	105	114	8.2	110	114	3.6	70 - 130	30	
Chlorobenzene	ND	1.0	97	103	6.0	102	101	1.0	70 - 130	30	

SDG I.D.: GCF10994

% % Blk **LCSD** LCS MSD **RPD** LCS MS MS Rec Blank RL % **RPD** % % RPD Limits Limits % Parameter Chloroethane ND 1.0 98 108 9.7 100 98 70 - 130 2.0 30 Chloroform ND 1.0 98 106 7.8 102 103 1.0 70 - 130 30 Chloromethane ND 1.0 96 104 8.0 88 91 3.4 40 - 160 30 0.40 ND 96 104 8.0 104 106 1.9 70 - 130 30 cis-1,3-Dichloropropene Dibromochloromethane ND 0.50 101 108 6.7 122 125 2.4 70 - 130 30 Dibromomethane ND 1.0 96 100 4.1 101 103 2.0 70 - 130 30 Dichlorodifluoromethane ND 1.0 110 117 6.2 96 91 5.3 40 - 160 30 ND 1.0 100 108 7.7 98 99 1.0 70 - 130 Ethyl ether 30 Ethylbenzene ND 97 7.0 105 103 70 - 130 30 1.0 104 1.9 Hexachlorobutadiene ND 0.40 97 109 11.7 110 112 1.8 70 - 130 30 Isopropylbenzene ND 1.0 96 105 9.0 107 106 0.9 70 - 130 30 ND 1.0 96 7.0 103 70 - 130 m&p-Xylene 103 102 1.0 30 Methyl ethyl ketone ND 5.0 88 97 9.7 90 98 8.5 40 - 160 30 Methyl t-butyl ether (MTBE) ND 1.0 93 99 6.3 95 101 6.1 70 - 130 30 Methylene chloride ND 1.0 90 96 6.5 91 93 22 70 - 130 30 Naphthalene ND 1.0 76 104 31.1 103 118 13.6 70 - 130 30 ND 95 10.9 n-Butylbenzene 1.0 106 108 109 0.9 70 - 130 30 n-Propylbenzene ND 1.0 97 105 7.9 105 104 1.0 70 - 130 30 ND 1.0 98 107 105 70 - 130 o-Xylene 8.8 106 0.9 30 p-Isopropyltoluene ND 1.0 96 105 9.0 107 106 0.9 70 - 130 30 sec-Butylbenzene ND 1.0 101 110 8.5 112 114 1.8 70 - 130 30 Styrene ND 1.0 98 104 5.9 102 103 1.0 70 - 130 30 ND 1.0 97 105 7.9 107 105 tert-Butylbenzene 1.9 70 - 130 30 Tetrahydrofuran (THF) ND 2.5 91 97 6.4 92 100 8.3 70 - 130 30 Toluene ND 1.0 98 105 6.9 111 108 2.7 70 - 130 30 trans-1,2-Dichloroethene ND 96 105 9.0 107 106 1.0 0.9 70 - 130 30 trans-1,3-Dichloropropene ND 0.40 95 102 7.1 72 72 0.0 70 - 130 30 97 trans-1,4-dichloro-2-butene ND 5.0 111 13.5 87 104 17.8 70 - 130 30 Trichlorofluoromethane ND 101 107 5.8 104 102 1.9 1.0 70 - 130 30 Trichlorotrifluoroethane ND 1.0 108 116 7.1 114 108 5.4 70 - 130 30 Vinyl chloride ND 1.0 92 101 9.3 92 90 2.2 70 - 130 30 % 1,2-dichlorobenzene-d4 94 100 102 103 % 102 2.0 1.0 70 - 130 30 % Bromofluorobenzene 94 % 99 100 1.0 99 100 1.0 70 - 130 30 % Dibromofluoromethane 103 103 % 104 1.0 100 106 5.8 70 - 130 30 % Toluene-d8 93 % 102 102 0.0 105 103 1.9 70 - 130 30 Comment: Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 514336 (ug/L), QC Sample No: CF12269 (CF10994 (10X))

Volatiles - Ground Wate	<u>er</u>						
cis-1,2-Dichloroethene	ND	1.0	90	97	7.5	70 - 130	30
Tetrachloroethene	ND	1.0	93	99	6.3	70 - 130	30
Trichloroethene	ND	1.0	91	96	5.3	70 - 130	30
% 1,2-dichlorobenzene-d4	95	%	99	101	2.0	70 - 130	30
% Bromofluorobenzene	93	%	94	96	2.1	70 - 130	30
% Dibromofluoromethane	106	%	99	101	2.0	70 - 130	30
% Toluene-d8	94	%	103	103	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%.

% % Blk LCS LCSD LCS MSD **RPD** MS MS Rec Blank RPD RL % % % % **RPD** Limits Limits Parameter QA/QC Batch 514152 (ug/L), QC Sample No: CF10994 (CF10994) Volatile Petroleum Hydrocarbons - Ground Water Unadjusted C5-C8 Aliphatics (*1) 100 100 103 3.0 95 98 70 - 130 20 3.1 Unadjusted C9-C12 Aliphatics (*1) ND 100 103 106 2.9 77 83 7.5 70 - 130 20 C5-C8 Aliphatic Hydrocarbons *1,2 ND 100 92 95 100 103 3.0 3.2 70 - 130 20 C9-C12 Aliphatic Hydrocarbons *1, ND 100 103 106 2.9 66 72 70 - 130 20 C9-C10 Aromatic Hydrocarbons *1 ND 100 96 99 93 94 70 - 130 3.1 1.1 20 ND 87 90 3.4 90 90 70 - 130 20 Benzene 1.0 0.0 Ethyl Benzene ND 1.0 93 95 2.1 96 95 70 - 130 20 1.0 ND 93 95 97 **MTBE** 1.0 2.1 95 2.1 70 - 130 20 Naphthalene ND 5.0 91 92 1.1 87 92 5.6 70 - 130 20 Toluene ND 1.0 93 95 2.1 96 95 1.0 70 - 130 20 m,p-Xylenes ND 2.0 94 97 3.1 96 96 0.0 70 - 130 20 ND 1.0 92 94 2.2 94 94 o-Xylene 0.0 70 - 130 20 % 2,5-Dibromotoluene (PID) 101 % 103 98 5.0 104 90 14.4 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

SDG I.D.: GCF10994

January 21, 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, January 21, 2020 Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report GCF10994 - STANTECMA

State: MA

State:	MA						RL	Analysis
SampNo	Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	Criteria	Units
CF10994	\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF10994	\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
CF10994	\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
CF10994	\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
CF10994	\$8260GWR	Vinyl chloride	MA / CMR 310.40.1600 / GW-1 (mg/l)	10	1.0	2	2	ug/L
CF10994	\$8260GWR	Trichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	67	10	5	5	ug/L
CF10994	\$8260GWR	Vinyl chloride	MA / CMR 310.40.1600 / GW-1 (mg/l)	10	1.0	2	2	ug/L
CF10994	\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
CF10994	\$8260GWR	Tetrachloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	100	10	5	5	ug/L
CF10994	\$8260GWR	cis-1,2-Dichloroethene	MA / CMR 310.40.1600 / GW-1 (mg/l)	24	10	20	20	ug/L
CF10994	\$8260GWR	Trichloroethene	MA / GROUNDWATER STANDARDS / GW-1	67	10	5	5	ug/L
CF10994	\$8260GWR	Vinyl chloride	MA / GROUNDWATER STANDARDS / GW-1	10	1.0	2	2	ug/L
CF10994	\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
CF10994	\$8260GWR	Tetrachloroethene	MA / GROUNDWATER STANDARDS / GW-1	100	10	5	5	ug/L
CF10994	\$MCPADD-W	M 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
CF10994	\$MCPADD-W	M 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.

	MassDEP Analytical Protocol Certification Form											
Laboratory Name: Phoenix Environmental Laboratories, Inc. Project #:												
Proje	ct Locati	i on: 30 PE	ENNIM	IAN		RTN:						
This F	orm provid	les certification	ons for	the following data set	: [list Laborato	ry San	nple ID Number(s)]				
CF109	94											
		undwater/Sur			t Drinking	Wate	r 🗌 Air	O	ther:			
	-	check all th		-	0004 B # 11		740011 0		5	ED 4011		
8260 V CAM II		7470/7471 H	g V	MassDEP VPH CAM IV A ✓	8081 Pesticides CAM V B		7196 Hex Cr CAM VI B		CAM IX	EP APH		
8270 S CAM II		7010 Metals CAM III C		MassDEP EPH CAM IV B ✓	8151 Herbicides CAM V C		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	е				
	Affirmati	ive respons	es to q	uestions A through	F are require	d for	"Presumptive	Certai	inty" s	tatus		
Α	Chain-of-0	Custody, pro	perly p	a condition consiste reserved (including to llyzed with method ho	emperature*) ir	the f	ield or	V	Yes	□ No		
В		analytical m CAM protoco		s) and all associated owed?	QC requiremer	nts sp	ecified in the	V	Yes	□No		
С		CAM protoco		actions and analytica plemented for all ider				✓,	Yes	□ No		
D	CAM VII A		ssuran	omply with all the repo ce and Quality Contro ata"?				✓ ,	Yes	□ No		
Ш		modification		ods only: Was each refer to the individual				✓ ,	Yes	□No		
	b. APH a method?	nd TO-15 me	ethods	only: Was the compl	ete analyte list	repor	ted for each		Yes	□ No		
F	conformal		ed and	tocol QC and perforn evaluated in a labora rough E)?				✓ ,	Yes	□ No		
	Resp	onses to q	uestio	ns G, H and I below	is required fo	r "Pre	sumptive Cert	tainty'	' statu	S		
G		reporting lim CAM protoco		r below all CAM repo	orting limits spe	cified	in the		Yes	✓ No		
				resumptive Certainty" cribed in 310 CMR 40.				data us	sability	and		
I	See Secti	ons: EPH, V	OA, VF	andards specified in t PH Narrations .	·				Yes	✓ No		
I	Were resu protocol(s)?		complete analyte lis	•				Yes	□ No		
I. the u	ındersiane		•	esponses must be addr ains and penalties of p					rv of th	iose		
respor	sible for o		nforma	tion, the material conf								
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Sign	ature: -	, 0	- V (W)	4 1 340 60	_		Project Manag					



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



MCP Certification Report

January 21, 2020 SDG I.D.: GCF10994

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.

EPH Narration

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

QC Batch 514708 (Samples: CF10994): -----

One or more analytes is below the method criteria. A low bias for these analytes is possible. (C36 - Hexatriacontane)

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. (C9 - Nonane)

Instrument:

AU-FID3 01/17/20-1

Adam Werner, Chemist 01/17/20

CF10994

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 01/17/20-1

Adam Werner, Chemist 01/17/20

CF10994

The initial calibration (ALN25BI) 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 514708 (CF10994)

CF10994

All LCS recoveries were within 40 - 140 with the following exceptions: C36 - Hexatriacontane(23%)

All LCSD recoveries were within 40 - 140 with the following exceptions: C36 - Hexatriacontane(25%), 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%

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:



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Certification Report

January 21, 2020 SDG I.D.: GCF10994

Mercury Narration

MERLIN 01/15/20 07:38

Rick Schweitzer. Chemist 01/15/20

CF10994

The method preparation blank contains all of the acids and reagents as the samples; the instrument blanks do not.

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 514110 (CF11603)

CF10994

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 01/14/20 08:53

Emily Kolominskaya, Chemist 01/14/20

CF10994

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 514163 (CF10467)

CF10994

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

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

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

ICPMS Metals Narration

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

Instrument:



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Certification Report

January 21, 2020 SDG I.D.: GCF10994

ICPMS Metals Narration

ICPMS 01/14/20 13:52

Cindy Pearce, Chemist 01/14/20

CF10994

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 513984 (CF10473)

CF10994

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

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

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

SVOASIM Narration

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

Instrument:

CHEM27 01/14/20-1

Wes Bryon, Chemist 01/14/20

CF10994

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_0103):

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 (CHEM27/0114_03-27_SIM18_0103) (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.

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

QC (Batch Specific):

Batch 513953 (CF10994)

CF10994

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

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

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%)



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MCP Certification Report

January 21, 2020 SDG I.D.: GCF10994

SVOASIM 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.

VOA Narration

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

QC Batch 514113 (Samples: CF10994): -----

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

Instrument:

CHEM17 01/13/20-2

Michael Hahn, Chemist 01/13/20

CF10994

Initial Calibration Evaluation (CHEM17/VT-S011220):

98% of target compounds met criteria.

The following compounds had %RSDs >20%: 1,2-Dibromo-3-chloropropane 25% (20%), trans-1,4-dichloro-2-butene 32% (20%) The following compounds did not meet recommended response factors: 1,2-Dibromo-3-chloropropane 0.035 (0.05), 2-Hexanone 0.087 (0.1), Acetone 0.052 (0.1), Bromoform 0.094 (0.1), Methyl ethyl ketone 0.078 (0.1), Tetrahydrofuran (THF) 0.046 (0.05), trans-1,4-dichloro-2-butene 0.041 (0.05)

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

Continuing Calibration Verification (CHEM17/0113_30-VT-S011220) (MCP Compliance):

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

95% 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: 1,2-Dibromo-3-chloropropane 0.032 (0.05), 2-Hexanone 0.077 (0.1), Acetone 0.058 (0.1), Bromoform 0.092 (0.1), Methyl ethyl ketone 0.070 (0.1), Tetrahydrofuran (THF) 0.043 (0.05), trans-1,4-dichloro-2-butene 0.031 (0.05)

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

CHEM17 01/14/20-2

Michael Hahn, Chemist 01/14/20

CF10994

Initial Calibration Evaluation (CHEM17/VT-S011220):

98% 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 (CHEM17/0114_31-VT-S011220) (MCP Compliance):

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

95% 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.

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



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MCP Certification Report

January 21, 2020 SDG I.D.: GCF10994

VOA Narration

QC (Batch Specific):

Batch 514113 (CF10770) CHEM17 1/13/2020-2

CF10994

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: Naphthalene(31.1%)

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

Batch 514336 (CF12269) CHEM17 1/14/2020-2

CF10994

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%.

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.

VPH Narration

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

QC Batch 514152 (Samples: CF10994): ----

The LCS/LCSD recovery is acceptable. One or more analytes in the site specific matrix spike recovery is below the method criteria, therefore a low bias is likely. (C9-C12 Aliphatic Hydrocarbons *1,3)

Instrument:

PIDFID 01/13/20-2

Raman Makol. Chemist 01/13/20

CF10994

QC (Batch Specific):

Batch 514152 (CF10994)

CF10994

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.

Cooler: Yes No	Temp 1° Pg of Data Delivery/Contact Options:	Phone:	Project P.O: [<i>¶</i> 56 <i>D18¶</i> 2	This section MUST be completed with	Bottle Quantities,	1/2/2017	Ridos / Janos	\$100 TU O	(a) (b)				MA Data Format	WRA eSMART		S-2 GW-1 ☐ S-2 GW-2 ☐ S-2 GW-3 ☐ Tier II Checklist S-3 GW-1 ☐ S-3 GW-2 ☐ S-3 GW-3 ☐ Full Data Package*	SW Protection City Other	000
	37	Email: info@phoenixlabs.com Fax (860) 645-0823 Client Services (860) 645-8726	Project: 30 Penim	Report to: Kichar Leanul Invoice to:	QUOTE#	Analysis Reminest	THE STATE OF THE S						Time: Ri CI	ure RCP Cert irial) GW Protection	Time:	GB Leachability	3 Days* CA-tww I/C DEC Objectives Objectives	
		Environmental Laboratories, Inc.	STORE	Address: Quing, M		Sampler's Signature Date:	de: Mater SE=Sediment SL=Sludge S=Soil SD=Solid Liquid X =(Other)	PHOENIX USE ONLY Customer Sample Sample Date SAMPLE# Identification Matrix Sampled Sa	GW 1/10				Relinquished by: Accepted by: Date:	III DELLA MANAGORIA	J VIII	Associated and the control of the co		



Thursday, January 23, 2020

Attn: Mr Richard Learned Stantec **400 Crown Colony Drive**

Suite 200

Quincy MA 02169

Project ID: **30 PENNIMAN** SDG ID: GCF10995

Sample ID#s: CF10995, CF11044

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



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SDG Comments

January 23, 2020

SDG I.D.: GCF10995

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.

Sample CF10995 was received past hold time for Chromium, Hexavalent (SM3500CRB).



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

Sample Id Cross Reference

January 23, 2020

SDG I.D.: GCF10995

Project ID: 30 PENNIMAN

Client Id	Lab Id	Matrix
D1	CF10995	GROUND WATER
TRIP BLANK	CF11044	WATER



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

Analysis Report

January 23, 2020

FOR: Attn: Mr Richard Learned

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample InformationCustody InformationDateTimeMatrix:GROUND WATERCollected by:01/10/2012:00Location Code:STANTECMAReceived by:CP01/13/2013:30

Rush Request: Standard Analyzed by: see "By" below

Laboratory Data

SDG ID: GCF10995

Phoenix ID: CF10995

Project ID: 30 PENNIMAN

195601892

Client ID: D1

P.O.#:

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	01/15/20	EK	SW6010D
Arsenic	0.011	0.004	mg/L	1	01/15/20	EK	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	01/15/20	EK	SW6010D
Chromium	0.007	0.001	mg/L	1	01/15/20	EK	SW6010D
Copper	0.005	0.005	mg/L	1	01/15/20	EK	SW6010D
Iron	10.0	0.010	mg/L	1	01/15/20	EK	SW6010D
Mercury	< 0.0002	0.0002	mg/L	1	01/15/20	RS	SW7470A
Nickel	0.011	0.001	mg/L	1	01/15/20	EK	SW6010D
Lead	0.009	0.002	mg/L	1	01/15/20	EK	SW6010D
Antimony	< 0.005	0.005	mg/L	1	01/15/20	EK	SW6010D
Selenium	< 0.010	0.010	mg/L	1	01/15/20	EK	SW6010D
Trivalent Chromium	0.007	0.001	mg/L	1	01/15/20		Calculation
Zinc	0.034	0.004	mg/L	1	01/15/20	EK	SW6010D
Chloride	291	15.0	mg/L	5	01/13/20	TB	SM4500CLE-11
Chlorine Residual	< 0.02	0.02	mg/L	1	01/13/20 18:33	0	SM4500CI-G-00
Chromium, Hexavalent	< 0.01	0.01	mg/L	1	01/13/20 18:17	0	SM3500CRB-11
Ammonia as Nitrogen	0.70	0.25	mg/L	5	01/16/20	KDB	E350.1
Total Cyanide	< 0.010	0.010	mg/L	1	01/16/20	O/GD	SW9010C/SW9012B
O&G, Non-polar Material	< 1.4	1.4	mg/L	1	01/16/20	MSF	E1664A
Total Suspended Solids	56	10	mg/L	2	01/14/20	ARG	SM 2540D-11
Mercury Digestion	Completed				01/15/20	Q/Q	SW7470A
PCB Extraction (LDL)	Completed				01/14/20	AT	SW3510C
Semi-Volatile Extraction	Completed				01/13/20	P/AK	SW3520C
Total Metals Digestion	Completed				01/14/20	AG	
Polychlorinated Biphe	enyls						
PCB-1016	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
PCB-1221	ND	0.048	ug/L	1	01/15/20	SC	SW8082A

Project ID: 30 PENNIMAN Phoenix I.D.: CF10995

Client ID: D1

Parameter			RL/					
PCB-1242	Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
PCB-1248	PCB-1232	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
PCB-1264 ND	PCB-1242	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
PCB-1260	PCB-1248	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
PCB-1282 ND	PCB-1254	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
PCB-128B	PCB-1260	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
SOCS Surrogate Society Socie	PCB-1262	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
96 DCBP (Surrogate Rec) 53 % 1 0/1/5/20 SC 30 - 150 % % DCBP (Surrogate Rec) (Confirmation) 56 % 1 0/1/15/20 SC 30 - 150 % % TCMX (Surrogate Rec) (Confirmation) 71 % 1 0/1/15/20 SC 30 - 150 % 1,2-Dibromoethane (EDB) ND 0.02 ug/L 1 0/1/15/20 SC 30 - 150 % Volatiles TILIA (Confirmation) ND 1.0 ug/L 1 0/1/15/20 MH SW8260C 1,1,1-2 Tetrachloroethane ND 1.0 ug/L 1 0/1/15/20 MH SW8260C 1,1,1-2 Trichloroethane ND 1.0 ug/L 1 0/1/15/20 MH SW8260C 1,1,1-2 Trichloroethane ND 1.0 ug/L 1 0/1/15/20 MH SW8260C 1,1,1-2 Trichloroethane ND 1.0 ug/L 1 0/1/15/20 MH SW8260C 1,1,2-3 Trichloroethane ND <td< td=""><td>PCB-1268</td><td>ND</td><td>0.048</td><td>ug/L</td><td>1</td><td>01/15/20</td><td>SC</td><td>SW8082A</td></td<>	PCB-1268	ND	0.048	ug/L	1	01/15/20	SC	SW8082A
Ko DGBP (Surrogate Rec) (Confirmation) 56 % 1 01/15/20 SC 30 - 150 % % TCMX (Surrogate Rec) (Confirmation) 71 % 1 01/15/20 SC 30 - 150 % *** TCMX (Surrogate Rec) (Confirmation) 71 % 1 01/15/20 SC 30 - 150 % *** TCMX (Surrogate Rec) (Confirmation) ND 0.02 ug/L 1 01/16/20 SC 30 - 150 % **** TCMX (Surrogate Rec) (Confirmation) ND 0.02 ug/L 1 01/16/20 SC 30 - 150 % **** TCMX (Surrogate Rec) (Confirmation) ND 0.02 ug/L 1 01/16/20 GG 5 SW8260 C **** TCMX (Surrogate Rec) (Confirmation) ND 1.0 ug/L 1 01/15/20 MH SW8260 C 1,1,1,2-Trichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260 C 1,1,2-Trichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260 C 1,1-Dichloroethane ND 1.0 ug/L 1 01/15/20 <td>QA/QC Surrogates</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	QA/QC Surrogates							
% TCMX (Surrogate Rec) 66 % 1 01/15/20 SC 30 - 150 % % TCMX (Surrogate Rec) (Confirmation) 71 % 1 01/15/20 SC 30 - 150 % L_2-Dibromoethane (EDB) ND 0.02 ug/L 1 01/16/20 MS Volatiles T.1,12-Tertachloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1,12-Trichloroethane ND 0.50 ug/L 1 01/15/20 MH SW8260C 1,1,2-Tetrachloroethane ND 0.50 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroptopane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,2-3-Trichloropropane ND 1.0 ug/L <td>% DCBP (Surrogate Rec)</td> <td>53</td> <td></td> <td>%</td> <td>1</td> <td>01/15/20</td> <td>SC</td> <td>30 - 150 %</td>	% DCBP (Surrogate Rec)	53		%	1	01/15/20	SC	30 - 150 %
1,2-Dibromoethane (EDB) ND 0.02 ug/L 1 01/15/20 SC 30-150 %	% DCBP (Surrogate Rec) (Confirmation)	56		%	1	01/15/20	SC	30 - 150 %
1,2-Dibromoethane (EDB)	% TCMX (Surrogate Rec)	66		%	1	01/15/20	SC	30 - 150 %
1,1,1,2-Tetrachloroethane	% TCMX (Surrogate Rec) (Confirmation)	71		%	1	01/15/20	SC	30 - 150 %
1,1,1,2-Tetrachloroethane								
1,1,1,2-Tetrachloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1,1-Trichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1,2-Trichloroethane ND 0.50 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroethane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,1-Dichloroethene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,2-3-Trichlorobropropene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,2,3-Trichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,2,3-Trichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,2,3-Trichlorobenzene ND 1.0 ug/L 1	1,2-Dibromoethane (EDB)	ND	0.02	ug/L	1	01/16/20	CG	SW8011
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1,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3,5-Trimethylbenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,4-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH	•			=	1			
1,3,5-Trimethylbenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,4-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 5.0 ug/L 1 01/15/20 MH <td< td=""><td></td><td></td><td></td><td></td><td>1</td><td></td><td></td><td></td></td<>					1			
1,3-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,3-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,4-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C	1,2-Dichloropropane				1			
1,3-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 1,4-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C <	1,3,5-Trimethylbenzene			ug/L	1		МН	
1,4-Dichlorobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20				=	1			
2,2-Dichloropropane ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C	1,3-Dichloropropane	ND	1.0	=	1	01/15/20	MH	SW8260C
2-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C <	1,4-Dichlorobenzene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
2-Hexanone ND 5.0 ug/L 1 01/15/20 MH SW8260C 2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	2,2-Dichloropropane	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
2-Isopropyltoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	2-Chlorotoluene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
4-Chlorotoluene ND 1.0 ug/L 1 01/15/20 MH SW8260C 4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C	2-Hexanone	ND	5.0	ug/L	1	01/15/20	MH	SW8260C
4-Methyl-2-pentanone ND 5.0 ug/L 1 01/15/20 MH SW8260C Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	2-Isopropyltoluene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Acetone ND 25 ug/L 1 01/15/20 MH SW8260C Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	4-Chlorotoluene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Acrylonitrile ND 1.0 ug/L 1 01/15/20 MH SW8260C Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	4-Methyl-2-pentanone	ND	5.0	ug/L	1	01/15/20	MH	SW8260C
Benzene ND 0.70 ug/L 1 01/15/20 MH SW8260C Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	Acetone	ND	25	ug/L	1	01/15/20	MH	SW8260C
Bromobenzene ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	Acrylonitrile	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Bromochloromethane ND 1.0 ug/L 1 01/15/20 MH SW8260C Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	Benzene	ND	0.70	ug/L	1	01/15/20	МН	SW8260C
Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	Bromobenzene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Bromodichloromethane ND 0.50 ug/L 1 01/15/20 MH SW8260C	Bromochloromethane	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
		ND	0.50		1	01/15/20	МН	SW8260C
	Bromoform	ND	1.0		1	01/15/20	МН	SW8260C
Bromomethane ND 1.0 ug/L 1 01/15/20 MH SW8260C	Bromomethane	ND	1.0	ug/L	1	01/15/20	МН	SW8260C

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		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Carbon Disulfide	ND	5.0	ug/L	1	01/15/20	MH	SW8260C
Carbon tetrachloride	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Chlorobenzene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Chloroethane	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Chloroform	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Chloromethane	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
cis-1,2-Dichloroethene	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
cis-1,3-Dichloropropene	ND	0.40	ug/L	1	01/15/20	MH	SW8260C
Dibromochloromethane	ND	0.50	ug/L	1	01/15/20	MH	SW8260C
Dibromomethane	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Dichlorodifluoromethane	ND	1.0	ug/L	1	01/15/20	MH	SW8260C
Ethylbenzene	1.0	1.0	ug/L	1	01/15/20	MH	SW8260C
Hexachlorobutadiene	ND	0.40	ug/L	1	01/15/20	MH	SW8260C
Isopropylbenzene	1.9	1.0	ug/L	1	01/15/20	MH	SW8260C
m&p-Xylene	2.8	1.0	ug/L	1	01/15/20	MH	SW8260C
Methyl ethyl ketone	ND	5.0	ug/L	1	01/15/20	MH	SW8260C
Methyl t-butyl ether (MTBE)	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Methylene chloride	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Naphthalene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
n-Butylbenzene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
n-Propylbenzene	1.7	1.0	ug/L	1	01/15/20	МН	SW8260C
o-Xylene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
p-Isopropyltoluene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
sec-Butylbenzene	1.0	1.0	ug/L	1	01/15/20	МН	SW8260C
Styrene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
tert-Butylbenzene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Tetrachloroethene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Tetrahydrofuran (THF)	ND	2.5	ug/L	1	01/15/20	МН	SW8260C
Toluene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Total Xylenes	2.8	1.0	ug/L	1	01/15/20	МН	SW8260C
trans-1,2-Dichloroethene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
trans-1,3-Dichloropropene	ND	0.40	ug/L	1	01/15/20	МН	SW8260C
trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	01/15/20	МН	SW8260C
Trichloroethene	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Trichlorofluoromethane	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Trichlorotrifluoroethane	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
Vinyl chloride	ND	1.0	ug/L	1	01/15/20	МН	SW8260C
QA/QC Surrogates			9				
% 1,2-dichlorobenzene-d4	97		%	1	01/15/20	МН	70 - 130 %
% Bromofluorobenzene	93		%	1	01/15/20	МН	70 - 130 %
% Dibromofluoromethane	104		%	1	01/15/20	МН	70 - 130 %
% Toluene-d8	96		%	1	01/15/20	МН	70 - 130 %
	-						
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	01/15/20	MH	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	01/15/20	МН	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	01/15/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	01/15/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	01/15/20	MH	SW8260C (OXY)

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Parameter	Result	RL/ PQL	Units Dilution		Date/Time	Ву	Reference
Ethanol	ND	400	ug/L	1	01/15/20	МН	SW8260C
Semivolatiles							
1,2,4,5-Tetrachlorobenzene	ND	3.3	ug/L	1	01/15/20	WB	SW8270D
1,2,4-Trichlorobenzene	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
1,2-Dichlorobenzene	ND	2.4	ug/L	1	01/15/20	WB	SW8270D
1,2-Diphenylhydrazine	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
1,3-Dichlorobenzene	ND	2.4	ug/L	1	01/15/20	WB	SW8270D
1,4-Dichlorobenzene	ND	2.4	ug/L	1	01/15/20	WB	SW8270D
2,4,5-Trichlorophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2,4,6-Trichlorophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2,4-Dichlorophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2,4-Dimethylphenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2,4-Dinitrophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2,4-Dinitrotoluene	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
2,6-Dinitrotoluene	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
2-Chloronaphthalene	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
2-Chlorophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2-Methylphenol (o-cresol)	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
2-Nitroaniline	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
2-Nitrophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
3&4-Methylphenol (m&p-cresol)	ND	9.5	ug/L	1	01/15/20	WB	SW8270D
3,3'-Dichlorobenzidine	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
3-Nitroaniline	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
4,6-Dinitro-2-methylphenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
4-Bromophenyl phenyl ether	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
4-Chloro-3-methylphenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
4-Chloroaniline	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
4-Chlorophenyl phenyl ether	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
4-Nitroaniline	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
4-Nitrophenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
Acetophenone	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Aniline	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Benzidine	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Benzoic acid	ND	48	ug/L	1	01/15/20	WB	SW8270D
Benzyl butyl phthalate	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Bis(2-chloroethoxy)methane	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Bis(2-chloroethyl)ether	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
Bis(2-chloroisopropyl)ether	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Bis(2-ethylhexyl)phthalate	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
Carbazole	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Dibenzofuran	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Diethyl phthalate	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Dimethylphthalate	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Di-n-butylphthalate	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Di-n-octylphthalate	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
Hexachloroethane	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
Isophorone	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
N-Nitrosodi-n-propylamine	ND	4.8	ug/L	1	01/15/20	WB	SW8270D
N-Nitrosodiphenylamine	ND	4.8	ug/L	1	01/15/20	WB	SW8270D

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	RL/						
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Pentachloronitrobenzene	ND	2.4	ug/L	1	01/15/20	WB	SW8270D
Phenol	ND	0.95	ug/L	1	01/15/20	WB	SW8270D
QA/QC Surrogates							
% 2,4,6-Tribromophenol	82		%	1	01/15/20	WB	15 - 110 %
% 2-Fluorobiphenyl	70		%	1	01/15/20	WB	30 - 130 %
% 2-Fluorophenol	46		%	1	01/15/20	WB	15 - 110 %
% Nitrobenzene-d5	60		%	1	01/15/20	WB	30 - 130 %
% Phenol-d5	52		%	1	01/15/20	WB	15 - 110 %
% Terphenyl-d14	71		%	1	01/15/20	WB	30 - 130 %
Semivolatiles (SIM)							
2-Methylnaphthalene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Acenaphthene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Acenaphthylene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Anthracene	ND	0.09	ug/L	1	01/15/20	PS	SW8270D (SIM)
Benz(a)anthracene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Benzo(a)pyrene	ND	0.19	ug/L	1	01/15/20	PS	SW8270D (SIM)
Benzo(b)fluoranthene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Benzo(ghi)perylene	ND	0.02	ug/L	1	01/15/20	PS	SW8270D (SIM)
Benzo(k)fluoranthene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Chrysene	ND	0.05	ug/L	1	01/15/20	PS	SW8270D (SIM)
Dibenz(a,h)anthracene	ND	0.02	ug/L	1	01/15/20	PS	SW8270D (SIM)
Fluoranthene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Fluorene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Hexachlorobenzene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Hexachlorobutadiene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Hexachlorocyclopentadiene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Indeno(1,2,3-cd)pyrene	ND	0.10	ug/L	1	01/15/20	PS	SW8270D (SIM)
Naphthalene	0.59	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Nitrobenzene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
N-Nitrosodimethylamine	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Pentachlorophenol	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Phenanthrene	ND	0.48	ug/L	1	01/15/20	PS	SW8270D (SIM)
Pyrene	ND	0.07	ug/L	1	01/15/20	PS	SW8270D (SIM)
Pyridine	ND	1.9	ug/L	1	01/15/20	PS	SW8270D (SIM)
QA/QC Surrogates							
% 2,4,6-Tribromophenol	79		%	1	01/15/20	PS	15 - 110 %
% 2-Fluorobiphenyl	67		%	1	01/15/20	PS	40 - 140 %
% 2-Fluorophenol	47		%	1	01/15/20	PS	15 - 110 %
% Nitrobenzene-d5	56		%	1	01/15/20	PS	40 - 140 %
% Phenol-d5	56		%	1	01/15/20	PS	15 - 110 %
% Terphenyl-d14	61		%	1	01/15/20	PS	40 - 140 %
1,4-dioxane							
1,4-dioxane	ND	0.20	ug/l	1	01/15/20	AW	SW8270DSIM
QA/QC Surrogates							
% 1,4-dioxane-d8	81		%	1	01/15/20	AW	30 - 130 %
Extraction for 1,4-Dioxane	Completed				01/14/20	S/S	

Project ID: 30 PENNIMAN Phoenix I.D.: CF10995

Client ID: D1

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

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.

8260 Analysis:

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

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

January 23, 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

January 23, 2020

FOR: Attn: Mr Joseph Salvetti

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample Information Custody Information Date <u>Time</u> WATER Collected by: 01/13/20 14:51 Matrix: Received by: Location Code: **STANTECMA** SW 01/13/20 13:30 Rush Request: Standard Analyzed by: see "By" below

Laboratory Data

SDG ID: GCF10995

Phoenix ID: CF11044

30 PENNIMAN Project ID: Client ID: TRIP BLANK

P.O.#:

RL/

Parameter	Result PQL Units Dilution		Dilution	Date/Time	Ву	Reference		
Volatiles								
1,1,1,2-Tetrachloroethane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,1,1-Trichloroethane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,1,2,2-Tetrachloroethane	ND	0.50	ug/L	1	01/13/20	МН	SW8260C	
1,1,2-Trichloroethane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,1-Dichloroethane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,1-Dichloroethene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,1-Dichloropropene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2,3-Trichlorobenzene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2,3-Trichloropropane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2,4-Trichlorobenzene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2,4-Trimethylbenzene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2-Dibromo-3-chloropropane	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
1,2-Dibromoethane	ND	0.50	ug/L	1	01/13/20	МН	SW8260C	
1,2-Dichlorobenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
1,2-Dichloroethane	ND	0.60	ug/L	1	01/13/20	MH	SW8260C	
1,2-Dichloropropane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
1,3,5-Trimethylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
1,3-Dichlorobenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
1,3-Dichloropropane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
1,4-Dichlorobenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
2,2-Dichloropropane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
2-Chlorotoluene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
2-Hexanone	ND	5.0	ug/L	1	01/13/20	МН	SW8260C	
2-Isopropyltoluene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C	
4-Chlorotoluene	ND	1.0	ug/L	1	01/13/20	МН	SW8260C	
4-Methyl-2-pentanone	ND	5.0	ug/L	1	01/13/20	МН	SW8260C	

Project ID: 30 PENNIMAN Client ID: TRIP BLANK

Acetone ND 25 ug/L 1 01/13/20 MH SW8280C Acrylonitrile ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromobenzene ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromodichioromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromodichioromethane ND 0.50 ug/L 1 01/13/20 MH SW8280C Bromodichioromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromodichioromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Carbon Istrachioride ND 1.0 ug/L 1 01/13/20 MH SW8280C Chlorochezane ND 1.0 ug/L 1 01/13/20 MH SW8280C Chlorochezane ND 1.0 ug/L 1 01/13/20 MH SW8280C </th <th>Parameter</th> <th>Result</th> <th>RL/ PQL</th> <th>Units</th> <th>Dilution</th> <th>Date/Time</th> <th>Ву</th> <th>Reference</th>	Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Accylonitrile	Acetone	ND	25	ug/L	1	01/13/20	МН	SW8260C
Benzene	Acrylonitrile	ND	1.0		1	01/13/20	МН	SW8260C
Bromochenzene ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromochloromethane ND 0.50 ug/L 1 01/13/20 MH SW8280C Bromochrom ND 1.0 ug/L 1 01/13/20 MH SW8280C Bromomethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Carbon letrachloride ND 1.0 ug/L 1 01/13/20 MH SW8280C Chlorobenzene ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroderm ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroderm ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroderhane ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroderhane ND 1.0 ug/L 1 01/13/20 MH SW8280C	-	ND	0.70		1	01/13/20	MH	SW8260C
Bromodichloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Bromodichloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C SW	Bromobenzene	ND	1.0		1	01/13/20	МН	SW8260C
Bromodorhane ND 0.50 ug/L 1 01/13/20 MH SW8280C Bromoform ND 1.0 ug/L 1 01/13/20 MH SW8280C Carbon Disulfide ND 5.0 ug/L 1 01/13/20 MH SW8280C Carbon Intrachloride ND 1.0 ug/L 1 01/13/20 MH SW8280C Chlorobenzene ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroethane ND 1.0 ug/L 1 01/13/20 MH SW8280C	Bromochloromethane	ND	1.0		1	01/13/20	MH	SW8260C
Bromoremethane	Bromodichloromethane	ND	0.50		1	01/13/20	MH	SW8260C
Carbon Disulfide ND 5.0 ug/L 1 01/13/20 MH SW8260C Carbon tetrachloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Cis-1,3-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Dibromochloromethane ND 0.40 ug/L 1 01/13/20 MH SW8260C Dibromochloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Elhylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Elhylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C	Bromoform	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Carbon tetrachloride	Bromomethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Chlorobenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Chloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C cis-1,3-Dichloroethene ND 0.40 ug/L 1 01/13/20 MH SW8260C Dibromochloromethane ND 0.40 ug/L 1 01/13/20 MH SW8260C Dibromochloromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C	Carbon Disulfide	ND	5.0	ug/L	1	01/13/20	MH	SW8260C
Chloroethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloroform ND 1.0 ug/L 1 01/13/20 MH SW8280C Chloromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C cis-1,2-Dichloroptropene ND 0.40 ug/L 1 01/13/20 MH SW8280C Dibromochloromethane ND 0.50 ug/L 1 01/13/20 MH SW8280C Dibromochloromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Dibromochlorodifluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8280C Eithylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8280C Hexachlorobutadiene ND 1.0 ug/L 1 01/13/20 MH SW8280C Hexachlorobutadiene ND 1.0 ug/L 1 01/13/20 MH <td>Carbon tetrachloride</td> <td>ND</td> <td>1.0</td> <td>ug/L</td> <td>1</td> <td>01/13/20</td> <td>MH</td> <td>SW8260C</td>	Carbon tetrachloride	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Chioroform	Chlorobenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Chloromethane	Chloroethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
cis-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C cis-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C Dibromoentlane ND 0.50 ug/L 1 01/13/20 MH SW8260C Dibromomethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl styl ketone ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl bethyl ketone ND 1.0 ug/L 1 01/13/20 MH SW	Chloroform	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
cis-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C Dibromochloromethane ND 0.50 ug/L 1 01/13/20 MH SW8260C Dichlorodiffuoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 M	Chloromethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Dibromochloromethane ND 0.50 ug/L 1 01/13/20 MH SW8260C Dibromomethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Dibromomethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Hexachlorobutadiene ND 0.40 ug/L 1 01/13/20 MH SW8260C Hexachlorobutadiene ND 1.0 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl teburg ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl teburg ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C N-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C N-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C N-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C N-Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C P-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrash-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01	cis-1,2-Dichloroethene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Dibromochloromethane ND 0.50 ug/L 1 01/13/20 MH SW8260C Dibromomethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Dibromomethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ethylbenzene ND 0.40 ug/L 1 01/13/20 MH SW8260C SW8260C Sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C MR.p-Xylene ND 1.0	cis-1,3-Dichloropropene	ND	0.40	ug/L	1	01/13/20	MH	SW8260C
Dichlorodifluoromethane		ND	0.50	ug/L	1	01/13/20	MH	SW8260C
Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Hexachlorobutadiene ND 0.40 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl tehyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20	Dibromomethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Ethylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Hexachlorobutadiene ND 0.40 ug/L 1 01/13/20 MH SW8260C Isopropylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl tehyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20	Dichlorodifluoromethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Isopropylbenzene		ND	1.0		1	01/13/20	MH	SW8260C
m&p-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 5.0 ug/L 1 01/13/20 MH SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C <td>•</td> <td>ND</td> <td>0.40</td> <td>ug/L</td> <td>1</td> <td>01/13/20</td> <td>MH</td> <td>SW8260C</td>	•	ND	0.40	ug/L	1	01/13/20	MH	SW8260C
m&p-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C Methyl ethyl ketone ND 5.0 ug/L 1 01/13/20 MH SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C O-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C	Isopropylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Methyl ethyl ketone ND 5.0 ug/L 1 01/13/20 MH SW8260C Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Ett-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C		ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Methyl t-butyl ether (MTBE) ND 1.0 ug/L 1 01/13/20 MH SW8260C Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C		ND	5.0	ug/L	1	01/13/20	MH	SW8260C
Methylene chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachlorotethene ND 1.0 ug/L 1 01/13/20 MH SW8260C	-	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Naphthalene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachlorotethene ND 1.0 ug/L 1 01/13/20 MH SW8260C		ND	1.0	ug/L	1	01/13/20	MH	SW8260C
n-Propylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrabutydrofuran (THF) ND 1.0 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C </td <td>Naphthalene</td> <td>ND</td> <td>1.0</td> <td>ug/L</td> <td>1</td> <td>01/13/20</td> <td>MH</td> <td>SW8260C</td>	Naphthalene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
o-Xylene ND 1.0 ug/L 1 01/13/20 MH SW8260C p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Totluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 0.40 ug/L 1 01/13/20 MH SW8260C	n-Butylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
p-Isopropyltoluene ND 1.0 ug/L 1 01/13/20 MH SW8260C sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Total will ene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-dichlorop-z-butene ND 5.0 ug/L 1 01/13/20 MH	n-Propylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
sec-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8	o-Xylene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Styrene ND 1.0 ug/L 1 01/13/20 MH SW8260C tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH <td>p-Isopropyltoluene</td> <td>ND</td> <td>1.0</td> <td>ug/L</td> <td>1</td> <td>01/13/20</td> <td>MH</td> <td>SW8260C</td>	p-Isopropyltoluene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
tert-Butylbenzene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20	sec-Butylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Tetrachloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroptenee ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloroptenee ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH	Styrene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Tetrahydrofuran (THF) ND 2.5 ug/L 1 01/13/20 MH SW8260C Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates W 1 01/13/20 MH 70 - 130 %	tert-Butylbenzene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Toluene ND 1.0 ug/L 1 01/13/20 MH SW8260C Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates S 1 01/13/20 MH 70 - 130 % <td>Tetrachloroethene</td> <td>ND</td> <td>1.0</td> <td>ug/L</td> <td>1</td> <td>01/13/20</td> <td>MH</td> <td>SW8260C</td>	Tetrachloroethene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Total Xylenes ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates 8 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Tetrahydrofuran (THF)	ND	2.5	ug/L	1	01/13/20	MH	SW8260C
trans-1,2-Dichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates V 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Toluene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
trans-1,3-Dichloropropene ND 0.40 ug/L 1 01/13/20 MH SW8260C trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates Surrogates W 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Total Xylenes	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
trans-1,4-dichloro-2-butene ND 5.0 ug/L 1 01/13/20 MH SW8260C Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates Surrogates W 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	trans-1,2-Dichloroethene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Trichloroethene ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates Surrogates W 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	trans-1,3-Dichloropropene	ND	0.40	ug/L	1	01/13/20	MH	SW8260C
Trichlorofluoromethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates SW8260C W 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	trans-1,4-dichloro-2-butene	ND	5.0	ug/L	1	01/13/20	MH	SW8260C
Trichlorotrifluoroethane ND 1.0 ug/L 1 01/13/20 MH SW8260C Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates Surrogates Valid In the control of the	Trichloroethene	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
Vinyl chloride ND 1.0 ug/L 1 01/13/20 MH SW8260C QA/QC Surrogates Surrogates Vinyl chlorobenzene-d4 94 % 1 01/13/20 MH 70 - 130 % 8 Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Trichlorofluoromethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
QA/QC Surrogates % 1,2-dichlorobenzene-d4 94 % 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Trichlorotrifluoroethane	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
% 1,2-dichlorobenzene-d4 94 % 1 01/13/20 MH 70 - 130 % % Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	Vinyl chloride	ND	1.0	ug/L	1	01/13/20	MH	SW8260C
% Bromofluorobenzene 92 % 1 01/13/20 MH 70 - 130 %	QA/QC Surrogates							
	% 1,2-dichlorobenzene-d4	94		%	1	01/13/20	MH	70 - 130 %
% Dibromofluoromethane 99 % 1 01/13/20 MH 70 - 130 %	% Bromofluorobenzene	92		%	1	01/13/20	MH	70 - 130 %
	% Dibromofluoromethane	99		%	1	01/13/20	МН	70 - 130 %

Project ID: 30 PENNIMAN

Client ID: TRIP BLANK

Phoenix I.D.: CF11044

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
% Toluene-d8	92		%	1	01/13/20	МН	70 - 130 %
Oxygenates & Dioxane							
1,4-Dioxane	ND	50	ug/L	1	01/13/20	МН	SW8260C (OXY)
Diethyl ether	ND	1.0	ug/L	1	01/13/20	MH	SW8260C (OXY)
Di-isopropyl ether	ND	1.0	ug/L	1	01/13/20	MH	SW8260C (OXY)
Ethyl tert-butyl ether	ND	1.0	ug/L	1	01/13/20	MH	SW8260C (OXY)
tert-amyl methyl ether	ND	1.0	ug/L	1	01/13/20	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:

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

January 23, 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

QA/QC Report

January 23, 2020

QA/QC Data

SDG I.D.: GCF10995

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 514110 (mg/L),	QC Sam	ole No: (CF11603	(CF1099	95)								
Mercury - Water	BRL	0.0002	<0.0002	<0.0002	NC	110			103			75 - 125	30
Comment:													
Additional Mercury criteria: LCS	acceptanc	e range f	or waters	is 80-120°	% and fo	or soils is	s 75-1259	%					
QA/QC Batch 514163 (mg/L),	QC Sam	ole No: (CF10467	(CF1099	95)								
ICP Metals - Aqueous													
Antimony	BRL	0.005	< 0.005	< 0.005	NC	102	104	1.9	114			75 - 125	20
Arsenic	BRL	0.004	< 0.004	< 0.004	NC	96.4	99.1	2.8	109			75 - 125	20
Cadmium	BRL	0.001	0.004	0.004	NC	97.7	101	3.3	90.6			75 - 125	20
Chromium	BRL	0.001	< 0.001	< 0.001	NC	97.9	101	3.1	94.5			75 - 125	20
Copper	BRL	0.005	< 0.005	< 0.005	NC	98.3	102	3.7	109			75 - 125	20
Iron	BRL	0.010	< 0.010	< 0.010	NC	104	104	0.0	95.7			75 - 125	20
Lead	BRL	0.002	0.003	0.003	NC	94.1	96.9	2.9	88.4			75 - 125	20
Nickel	BRL	0.001	0.100	0.100	0	100	103	3.0	91.4			75 - 125	20
Selenium	BRL	0.010	0.012	0.012	NC	94.0	96.1	2.2	101			75 - 125	20
Silver	BRL	0.001	< 0.001	0.002	NC	95.0	97.8	2.9	112			75 - 125	20
Zinc	BRL	0.004	0.292	0.292	0	96.9	100	3.1	104			75 - 125	20



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QA/QC Report

January 23, 2020

QA/QC Data

SDG I.D.: GCF10995

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 514435 (mg/L), Q	C Sam	ole No: (CF10215	(CF109	95)								
Total Cyanide Comment:	BRL	0.010	<0.010	<0.010	NC	90.4			97.5			90 - 110	30
Additional soil criteria LCS accepta	nce ran	ge is 80-	120% MS	acceptan	ce range	75-125	5%.						
QA/QC Batch 514080 (mg/L), Q	C Sam	ole No: (CF10779	(CF1099	95)								
Total Suspended Solids	BRL	2.5	17	14	NC	104						85 - 115	
QA/QC Batch 514509 (mg/L), Q	C Sam	ole No: (CF11778	(CF1099	95)								
O&G, Non-polar Material Comment:	BRL	1.4	<1.4	<1.4	NC	93.0			89.0			85 - 115	20
Additional: LCS acceptance range	is 85-11	5% MS a	acceptance	e range 7	5-125%								
QA/QC Batch 513988 (mg/L), Q	C Sam	ole No: (CF11334	(CF1099	95)								
Chromium, Hexavalent Comment:	BRL	0.01	<0.01	<0.01	NC	103			109			90 - 110	30
Additional Hexavalent Chromium of	riteria: L	CS acce	ptance ran	ige for wa	ters is 9	0-110%	and MS	accepta	nce ranç	ge is 85-	115%.		
QA/QC Batch 514023 (mg/L), Q	C Sam	ole No: (CF10330	(CF1099	95)								
Chloride	BRL	3.0	19.3	19.9	3.10	102			105			90 - 110	20
QA/QC Batch 514457 (mg/L), Q	C Sam	ole No: (CF10393	(CF1099	95)								
Ammonia as Nitrogen	BRL	0.05	<0.10	<0.10	NC	101			98.0			90 - 110	20
QA/QC Batch 514000 (mg/L), Q	C Sam	ole No: (CF11334	(CF1099	95)								
Chlorine Residual	BRL	0.02	< 0.02	< 0.02	NC	97.1							



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QA/QC Report

January 23, 2020

QA/QC Data

SDG I.D.: GCF10995 LCS **LCSD** LCS MS MSD MS Rec **RPD** Blank **RPD RPD** Limits RΙ % % % Limits Parameter % QA/QC Batch 514515 (ug/L), QC Sample No: CF13729 (CF10995) EDB and DBCP Analysis - Ground Water 1,2-Dibromoethane (EDB) 98 96 2.1 103 106 2.9 70 - 130 25 QA/QC Batch 514154 (ug/L), QC Sample No: CF10995 (CF10995) Polychlorinated Biphenyls - Ground Water PCB-1016 ND 0.050 99 98 1.0 40 - 140 20 PCB-1221 ND 0.050 40 - 140 20 PCB-1232 ND 0.050 40 - 140 20 PCB-1242 ND 0.050 40 - 140 20 ND 0.050 PCB-1248 40 - 140 20 PCB-1254 ND 0.050 40 - 140 20 PCB-1260 ND 0.050 98 110 11.5 40 - 140 20 PCB-1262 ND 0.050 40 - 140 20 PCB-1268 ND 0.050 40 - 140 20 % DCBP (Surrogate Rec) 71 % 96 101 5.1 30 - 150 20 % DCBP (Surrogate Rec) (Confirm 72 % 98 106 7.8 30 - 150 20 % 91 94 % TCMX (Surrogate Rec) 64 3.2 30 - 150 20 % TCMX (Surrogate Rec) (Confirm 64 % 94 104 10.1 30 - 150 20 Comment: A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate. QA/QC Batch 513952 (ug/L), QC Sample No: CF10667 (CF10995) Semivolatiles - Ground Water 1,2,4,5-Tetrachlorobenzene 3.5 78 40 - 140 20 ND 3.5 1,2,4-Trichlorobenzene 65 40 - 140 20 1,2-Dichlorobenzene ND 1.0 53 40 - 140 20 1,2-Diphenylhydrazine ND 1.6 75 40 - 140 20 1,3-Dichlorobenzene ND 1.0 49 40 - 140 20 1,4-Dichlorobenzene ND 1.0 50 40 - 140 20 ND 94 30 - 130 2,4,5-Trichlorophenol 1.0 20 2,4,6-Trichlorophenol ND 1.0 91 30 - 130 20 ND 2,4-Dichlorophenol 1.0 82 30 - 130 20 2,4-Dimethylphenol ND 85 1.0 30 - 130 20 2,4-Dinitrophenol ND 1.0 71 30 - 130 20 2,4-Dinitrotoluene ND 3.5 105 40 - 140 20 ND 3.5 102 2,6-Dinitrotoluene 40 - 140 20 2-Chloronaphthalene ND 3.5 71 40 - 140 20 2-Chlorophenol ND 1.0 59 30 - 130 20 2-Methylphenol (o-cresol) ND 1.0 75 30 - 13020 2-Nitroaniline ND 3.5 107 40 - 140 20 71 2-Nitrophenol ND 1.0 30 - 130 20 3&4-Methylphenol (m&p-cresol) ND 1.0 78 30 - 130 20 3,3'-Dichlorobenzidine 79 ND 5.0 40 - 140 20

SDG I.D.: GCF10995

%

% Blk **LCSD RPD** LCS LCS MS **MSD** MS Rec Blank RL % **RPD** % % RPD Limits Limits % Parameter 3-Nitroaniline ND 5.0 106 40 - 140 20 4,6-Dinitro-2-methylphenol ND 1.0 78 30 - 130 20 4-Bromophenyl phenyl ether ND 3.5 94 40 - 140 20 ND 1.0 96 30 - 130 20 4-Chloro-3-methylphenol 4-Chloroaniline ND 3.5 77 40 - 140 20 ND 1.0 85 40 - 140 20 4-Chlorophenyl phenyl ether 4-Nitroaniline ND 5.0 90 40 - 140 20 ND 1.0 98 4-Nitrophenol 30 - 130 20 ND 40 - 140 20 Acetophenone 3.5 67 Aniline ND 3.5 62 40 - 140 20 Benzidine ND 4.5 104 40 - 140 20 ND 30 - 130 Benzoic acid 10 52 20 Benzyl butyl phthalate ND 1.5 90 40 - 140 20 Bis(2-chloroethoxy)methane ND 3.5 68 40 - 140 20 Bis(2-chloroethyl)ether ND 1.0 48 40 - 140 20 Bis(2-chloroisopropyl)ether ND 1.0 50 40 - 140 20 ND 93 Bis(2-ethylhexyl)phthalate 1.5 40 - 140 20 ND 5.0 87 Carbazole 40 - 140 20 ND Dibenzofuran 3.5 76 40 - 140 20 Diethyl phthalate ND 1.5 87 40 - 140 20 ND 1.5 88 Dimethylphthalate 40 - 140 20 ND Di-n-butylphthalate 1.5 94 20 40 - 140 ND 1.5 99 Di-n-octylphthalate 40 - 140 20 Hexachloroethane ND 3.5 49 40 - 140 20 Isophorone ND 3.5 72 40 - 140 20 ND N-Nitrosodi-n-propylamine 3.5 76 40 - 140 20 N-Nitrosodiphenylamine ND 3.5 85 40 - 140 20 ND 5.0 96 Pentachloronitrobenzene 20 40 - 140 ND 1.0 51 Phenol 30 - 130 20 % 2,4,6-Tribromophenol 75 % 82 15 - 110 20 % 2-Fluorobiphenyl 60 % 66 30 - 130 20 % 2-Fluorophenol 33 % 39 15 - 110 20 % Nitrobenzene-d5 45 % 63 30 - 130 20 19 % Phenol-d5 % 49 15 - 110 20 % Terphenyl-d14 77 % 83 30 - 130 20 Comment: This batch consists of a Blank and LCS 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 514189 (ug/l), QC Sample No: CF12265 (CF10995) 1,4dioxane - Ground Water 1,4-dioxane ND 0.20 83 82 1.2 30 - 130 20 % 1,4-dioxane-d8 83 % 83 81 2.4 30 - 130 20 QA/QC Batch 513952 (ug/L), QC Sample No: CF10667 (CF10995) Semivolatiles (SIM) - Ground Water 2-Methylnaphthalene ND 78 40 - 140 20 0.50 Acenaphthene ND 0.50 81 40 - 140 20 ND 79 Acenaphthylene 0.50 40 - 140 20 Anthracene ND 0.50 96 40 - 140 20 Benz(a)anthracene ND 0.50 97 40 - 140 20

94

Benzo(a)pyrene

ND

0.50

40 - 140

20

SDG I.D.: GCF10995

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
Benzo(b)fluoranthene	ND	0.50	94						40 - 140	20	
Benzo(ghi)perylene	ND	0.50	94						40 - 140	20	
Benzo(k)fluoranthene	ND	0.50	97						40 - 140	20	
Chrysene	ND	0.50	96						40 - 140	20	
Dibenz(a,h)anthracene	ND	0.50	99						40 - 140	20	
Fluoranthene	ND	0.50	92						40 - 140	20	
Fluorene	ND	0.50	85						40 - 140	20	
Hexachlorobenzene	ND	0.50	87						40 - 140	20	
Hexachlorobutadiene	ND	0.50	66						40 - 140	20	
Hexachlorocyclopentadiene	ND	0.50	33						40 - 140	20	1
Indeno(1,2,3-cd)pyrene	ND	0.50	87						40 - 140	20	
Naphthalene	ND	0.50	73						40 - 140	20	
Nitrobenzene	ND	0.50	57						40 - 140	20	
N-Nitrosodimethylamine	ND	0.05	64						40 - 140	20	
Pentachlorophenol	ND	0.50	112						40 - 140	20	
Phenanthrene	ND	0.50	94						40 - 140	20	
Pyrene	ND	0.50	95						40 - 140	20	
Pyridine	ND	0.50	45						40 - 140	20	
% 2,4,6-Tribromophenol	61	%	82						15 - 110	20	
% 2-Fluorobiphenyl	55	%	73						40 - 140	20	
% 2-Fluorophenol	32	%	41						15 - 110	20	
% Nitrobenzene-d5	35	%	52						40 - 140	20	s
% Phenol-d5	18	%	52						15 - 110	20	
% Terphenyl-d14	64	%	81						40 - 140	20	
Comment:											

This batch consists of a Blank and LCS

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 514113 (ug/L), QC Sample No: CF10770 (CF11044)

Volatiles - Water											
1,1,1,2-Tetrachloroethane	ND	1.0	101	107	5.8	106	106	0.0	70 - 130	30	
1,1,1-Trichloroethane	ND	1.0	96	105	9.0	105	105	0.0	70 - 130	30	
1,1,2,2-Tetrachloroethane	ND	0.50	100	108	7.7	105	105	0.0	70 - 130	30	
1,1,2-Trichloroethane	ND	1.0	97	104	7.0	94	96	2.1	70 - 130	30	
1,1-Dichloroethane	ND	1.0	97	104	7.0	102	102	0.0	70 - 130	30	
1,1-Dichloroethene	ND	1.0	101	111	9.4	112	112	0.0	70 - 130	30	
1,1-Dichloropropene	ND	1.0	95	102	7.1	112	108	3.6	70 - 130	30	
1,2,3-Trichlorobenzene	ND	1.0	78	101	25.7	88	108	20.4	70 - 130	30	
1,2,3-Trichloropropane	ND	1.0	91	95	4.3	90	95	5.4	70 - 130	30	
1,2,4-Trichlorobenzene	ND	1.0	86	104	18.9	97	109	11.7	70 - 130	30	
1,2,4-Trimethylbenzene	ND	1.0	96	104	8.0	103	100	3.0	70 - 130	30	
1,2-Dibromo-3-chloropropane	ND	1.0	93	110	16.7	101	114	12.1	70 - 130	30	
1,2-Dibromoethane	ND	1.0	99	103	4.0	100	102	2.0	70 - 130	30	
1,2-Dichlorobenzene	ND	1.0	96	103	7.0	100	101	1.0	70 - 130	30	
1,2-Dichloroethane	ND	1.0	98	104	5.9	101	102	1.0	70 - 130	30	
1,2-Dichloropropane	ND	1.0	96	103	7.0	107	106	0.9	70 - 130	30	
1,3,5-Trimethylbenzene	ND	1.0	96	104	8.0	104	102	1.9	70 - 130	30	
1,3-Dichlorobenzene	ND	1.0	95	101	6.1	100	100	0.0	70 - 130	30	
1,3-Dichloropropane	ND	1.0	97	103	6.0	96	97	1.0	70 - 130	30	
1,4-Dichlorobenzene	ND	1.0	93	103	10.2	100	101	1.0	70 - 130	30	
1,4-dioxane	ND	100	106	114	7.3	105	101	3.9	40 - 160	30	
2,2-Dichloropropane	ND	1.0	90	100	10.5	90	90	0.0	70 - 130	30	

SDG I.D.: GCF10995

% % Blk **LCSD** RPD LCS LCS MS MSD MS Rec Blank RL **RPD** % % RPD Limits Limits % % Parameter ND 1.0 98 105 104 104 2-Chlorotoluene 6.9 0.0 70 - 130 30 ND 5.0 85 95 11.1 89 94 5.5 40 - 160 30 2-Hexanone 2-Isopropyltoluene ND 1.0 99 108 8.7 108 107 0.9 70 - 130 30 4-Chlorotoluene ND 1.0 93 101 100 99 70 - 130 30 8.2 1 0 4-Methyl-2-pentanone ND 5.0 89 97 8.6 46 47 2.2 40 - 160 30 ND 5.0 105 96 9.0 93 107 14.0 40 - 160 30 Acetone ND 5.0 99 106 99 107 7.8 70 - 130 30 Acrylonitrile 6.8 ND 0.70 95 7.1 106 Benzene 102 108 1.9 70 - 130 30 99 4.9 70 - 130 Bromobenzene ND 1.0 104 102 103 1.0 30 Bromochloromethane ND 1.0 99 108 8.7 103 102 1.0 70 - 130 30 Bromodichloromethane ND 0.50 98 106 7.8 106 106 0.0 70 - 130 30 ND 97 98 70 - 130 **Bromoform** 1.0 103 6.0 100 2.0 30 Bromomethane ND 1.0 100 113 12.2 80 87 8.4 40 - 160 30 Carbon Disulfide ND 1.0 92 101 9.3 99 101 2.0 70 - 130 30 1.0 Carbon tetrachloride ND 105 114 8.2 110 114 3.6 70 - 130 30 Chlorobenzene ND 1.0 97 103 6.0 102 101 1.0 70 - 130 30 ND Chloroethane 1.0 98 108 9.7 100 98 2.0 70 - 130 30 Chloroform ND 98 106 102 103 1.0 7.8 1.0 70 - 130 30 96 ND 104 91 Chloromethane 1.0 8.0 88 3.4 40 - 160 30 cis-1,2-Dichloroethene ND 1.0 97 104 7.0 NC NC NC 70 - 130 30 cis-1,3-Dichloropropene ND 0.40 96 104 104 106 8.0 1.9 70 - 130 30 Dibromochloromethane ND 0.50 101 108 6.7 122 125 2.4 70 - 130 30 103 Dibromomethane ND 1.0 96 100 4.1 101 2.0 70 - 130 30 Dichlorodifluoromethane ND 1.0 110 117 6.2 96 91 5.3 40 - 160 30 Ethyl ether ND 1.0 100 108 7.7 98 99 1.0 70 - 130 30 ND 97 104 Ethylbenzene 1.0 7.0 105 103 1.9 70 - 130 30 Hexachlorobutadiene ND 0.40 97 109 11.7 110 70 - 130 112 1.8 30 ND 96 105 9.0 107 106 Isopropylbenzene 1.0 0.9 70 - 130 30 ND 96 103 102 m&p-Xylene 1.0 103 7.0 1.0 70 - 130 30 Methyl ethyl ketone ND 5.0 88 97 9.7 90 98 8.5 40 - 160 30 99 Methyl t-butyl ether (MTBE) ND 1.0 93 6.3 95 101 6.1 70 - 130 30 Methylene chloride ND 90 96 91 1.0 6.5 93 2.2 70 - 130 30 Naphthalene ND 1.0 76 104 31.1 103 118 13.6 70 - 130 30 ND 95 n-Butylbenzene 1.0 106 10.9 108 109 0.9 70 - 130 30 n-Propylbenzene ND 1.0 97 105 7.9 105 104 1.0 70 - 130 30 o-Xylene ND 1.0 98 107 8.8 106 105 0.9 70 - 130 30 ND 9.0 106 p-Isopropyltoluene 1.0 96 105 107 0.9 70 - 130 30 sec-Butylbenzene ND 1.0 101 110 8.5 114 112 1.8 70 - 130 30 ND 98 104 5.9 102 70 - 130 Styrene 1.0 103 1.0 30 tert-Butylbenzene ND 1.0 97 105 7.9 107 105 1.9 70 - 130 30 ND 97 7.0 NC Tetrachloroethene 1.0 104 NC NC 70 - 130 30 91 Tetrahvdrofuran (THF) ND 2.5 97 6.4 92 100 8.3 70 - 130 30 Toluene ND 1.0 98 105 6.9 108 111 2.7 70 - 130 30 trans-1,2-Dichloroethene ND 1.0 96 105 9.0 107 106 0.9 70 - 130 30 trans-1,3-Dichloropropene ND 0.40 95 102 7.1 72 72 0.0 70 - 130 30 trans-1,4-dichloro-2-butene ND 5.0 97 111 13.5 87 104 17.8 70 - 130 30 Trichloroethene ND 1.0 94 102 8.2 108 104 3.8 70 - 130 30 107 ND 101 Trichlorofluoromethane 1.0 5.8 104 102 1.9 70 - 130 30 Trichlorotrifluoroethane ND 1.0 108 7.1 114 108 116 5.4 70 - 130 30 Vinyl chloride ND 1.0 92 101 9.3 92 90 2.2 70 - 130 30 94 % % 1,2-dichlorobenzene-d4 102 100 2.0 102 103 1.0 70 - 130 30 % Bromofluorobenzene 94 % 99 99 100 100 1.0 1.0 70 - 130 30 % Dibromofluoromethane 103 % 103 106 70 - 130 30 104 1 0 100 5.8

SDG I.D.: GCF10995

% % Blk **LCSD** LCS **RPD** LCS MS MSD MS Rec % Blank RL **RPD** % % RPD Limits % Limits Parameter % Toluene-d8 93 % 102 102 105 103 70 - 130 0.0 1.9 30 Comment: Additional 8260 criteria: 10% of compounds can be outside of acceptance criteria as long as recovery is 10%. QA/QC Batch 514336 (ug/L), QC Sample No: CF12269 (CF10995) Volatiles - Ground Water ND 1.0 97 7.5 1.1.1.2-Tetrachloroethane 90 70 - 13030 1.1.1-Trichloroethane ND 1.0 88 93 5.5 70 - 130 30 ND 0.50 1,1,2,2-Tetrachloroethane 91 103 12.4 70 - 130 30 1,1,2-Trichloroethane ND 1.0 82 100 19.8 70 - 130 30 1,1-Dichloroethane ND 1.0 88 95 7.7 70 - 130 30 ND 103 106 2.9 1.1-Dichloroethene 1.0 70 - 130 30 1,1-Dichloropropene ND 1.0 93 96 3.2 70 - 130 30 ND 1.0 82 95 14.7 1,2,3-Trichlorobenzene 70 - 130 30 1,2,3-Trichloropropane ND 1.0 86 88 2.3 70 - 130 30 ND 85 98 1,2,4-Trichlorobenzene 1.0 14.2 70 - 130 30 95 1,2,4-Trimethylbenzene ND 1.0 92 3.2 70 - 130 30 1,2-Dibromo-3-chloropropane ND 1.0 96 12.7 109 70 - 130 30 1.2-Dibromoethane ND 1.0 87 98 11.9 70 - 130 30 1,2-Dichlorobenzene ND 1.0 88 95 7.7 70 - 130 30 1.2-Dichloroethane ND 1.0 82 93 12.6 70 - 130 30 1,2-Dichloropropane ND 1.0 89 97 8.6 70 - 130 30 ND 94 95 1,3,5-Trimethylbenzene 1.0 1.1 70 - 130 30 1,3-Dichlorobenzene ND 1.0 89 94 5.5 70 - 130 30 1,3-Dichloropropane ND 1.0 85 95 11.1 70 - 130 30 1,4-Dichlorobenzene ND 1.0 88 94 6.6 70 - 130 30 1,4-dioxane ND 100 97 108 10.7 40 - 160 30 ND 1.0 92 70 - 130 30 2,2-Dichloropropane 86 6.7 2-Chlorotoluene ND 1.0 97 98 1.0 70 - 130 30 ND 5.0 73 90 20.9 40 - 160 30 2-Hexanone ND 1.0 95 98 3.1 70 - 130 30 2-Isopropyltoluene ND 1.0 90 93 70 - 130 4-Chlorotoluene 3.3 30 4-Methyl-2-pentanone ND 73 92 40 - 160 30 5.0 23.0 Acetone ND 5.0 83 92 40 - 160 30 10.3 Acrylonitrile ND 5.0 85 101 17.2 70 - 130 30 ND 92 97 70 - 130 Benzene 0.70 5.3 30 Bromobenzene ND 1.0 92 97 5.3 70 - 130 30 Bromochloromethane ND 1.0 85 97 13.2 70 - 130 30 Bromodichloromethane ND 0.50 86 98 13.0 70 - 130 30 **Bromoform** ND 1.0 83 95 13.5 70 - 130 30 Bromomethane ND 1.0 112 115 40 - 160 30 2.6 Carbon Disulfide ND 1.0 96 99 3.1 70 - 130 30 ND 94 1.0 101 7.2 Carbon tetrachloride 70 - 130 30 Chlorobenzene ND 1.0 91 95 4.3 70 - 130 30 Chloroethane ND 93 1.0 98 5.2 70 - 130 30 Chloroform ND 1.0 78 96 20.7 70 - 130 30 ND 1.0 87 94 7.7 Chloromethane 40 - 160 30 cis-1,2-Dichloroethene ND 1.0 90 97 7.5 70 - 130 30 cis-1,3-Dichloropropene ND 0.40 88 99 11.8 70 - 130 30 Dibromochloromethane ND 0.50 88 99 11.8 70 - 130 30 70 - 130 Dibromomethane ND 1.0 85 95 11.1 30 Dichlorodifluoromethane ND 1.0 101 106 4.8 40 - 160 30 82 Ethyl ether ND 1.0 96 15.7 70 - 130 30

SDG I.D.: GCF10995

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits
Ethylbenzene	ND	1.0	93	95	2.1				70 - 130	30
Hexachlorobutadiene	ND	0.40	110	103	6.6				70 - 130	30
Isopropylbenzene	ND	1.0	96	99	3.1				70 - 130	30
m&p-Xylene	ND	1.0	91	94	3.2				70 - 130	30
Methyl ethyl ketone	ND	5.0	67	86	24.8				40 - 160	30
Methyl t-butyl ether (MTBE)	ND	1.0	76	92	19.0				70 - 130	30
Methylene chloride	ND	1.0	83	91	9.2				70 - 130	30
Naphthalene	ND	1.0	89	103	14.6				70 - 130	30
n-Butylbenzene	ND	1.0	94	97	3.1				70 - 130	30
n-Propylbenzene	ND	1.0	95	96	1.0				70 - 130	30
o-Xylene	ND	1.0	93	96	3.2				70 - 130	30
p-Isopropyltoluene	ND	1.0	94	97	3.1				70 - 130	30
sec-Butylbenzene	ND	1.0	98	101	3.0				70 - 130	30
Styrene	ND	1.0	89	95	6.5				70 - 130	30
tert-Butylbenzene	ND	1.0	93	95	2.1				70 - 130	30
Tetrachloroethene	ND	1.0	93	99	6.3				70 - 130	30
Tetrahydrofuran (THF)	ND	2.5	71	87	20.3				70 - 130	30
Toluene	ND	1.0	94	99	5.2				70 - 130	30
trans-1,2-Dichloroethene	ND	1.0	93	98	5.2				70 - 130	30
trans-1,3-Dichloropropene	ND	0.40	85	99	15.2				70 - 130	30
trans-1,4-dichloro-2-butene	ND	5.0	97	127	26.8				70 - 130	30
Trichloroethene	ND	1.0	91	96	5.3				70 - 130	30
Trichlorofluoromethane	ND	1.0	97	101	4.0				70 - 130	30
Trichlorotrifluoroethane	ND	1.0	100	108	7.7				70 - 130	30
Vinyl chloride	ND	1.0	104	109	4.7				70 - 130	30
% 1,2-dichlorobenzene-d4	95	%	99	101	2.0				70 - 130	30
% Bromofluorobenzene	93	%	94	96	2.1				70 - 130	30
% Dibromofluoromethane	106	%	99	101	2.0				70 - 130	30
% Toluene-d8	94	%	103	103	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 514382 (ug/L), QC Sample No: CF10995 (CF10995)

Oxygenates - Ground Water

Ethanol 106 115 8.1 89 20.2 70 - 130 30

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

January 23, 2020

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

s = This parameter is outside laboratory Blank Surrogate specified recovery limits.

Thursday, January 23, 2020 Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report

State: MA

Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	RL Criteria	Analysis Units
\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
\$8270-SIMFSR	Benzoic acid	MA / CAM Protocol / SVOA AQ RL	ND	48		10	ug/L
AS-WM	Arsenic	MA / CMR 310.40.1600 / GW-1 (mg/l)	0.011	0.004	0.01	0.01	mg/L
AS-WM	Arsenic	MA / GROUNDWATER STANDARDS / GW-1	0.011	0.004	0.01	0.01	mg/L
\$8260GWR	Acetone	MA / CAM Protocol / VOA AQ RL	ND	25		10	ug/L
\$8260GWR	Carbon Disulfide	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
\$8260GWR	Tetrahydrofuran (THF)	MA / CAM Protocol / VOA AQ RL	ND	2.5		2	ug/L
\$8260GWR	trans-1,4-dichloro-2-butene	MA / CAM Protocol / VOA AQ RL	ND	5.0		2	ug/L
\$8260GWR	1,2-Dibromoethane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	0.50	0.02	0.02	ug/L
\$8260GWR	1,2-Dibromoethane	MA / GROUNDWATER STANDARDS / GW-1	ND	0.50	0.02	0.02	ug/L
\$MCPADD-WM	1 1,4-Dioxane	MA / CMR 310.40.1600 / GW-1 (mg/l)	ND	50	3	3	ug/L
\$MCPADD-WM	1 1,4-Dioxane	MA / GROUNDWATER STANDARDS / GW-1	ND	50	0.3	0.3	ug/L
	\$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8270-SIMFSR AS-WM AS-WM \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR \$8260GWR	\$8260GWR Acetone \$8260GWR Carbon Disulfide \$8260GWR Tetrahydrofuran (THF) \$8260GWR trans-1,4-dichloro-2-butene \$8260GWR 1,2-Dibromoethane \$8260GWR 1,2-Dibromoethane \$8270-SIMFSR Benzoic acid AS-WM Arsenic AS-WM Arsenic \$8260GWR Carbon Disulfide \$8260GWR Tetrahydrofuran (THF) \$8260GWR trans-1,4-dichloro-2-butene \$8260GWR 1,2-Dibromoethane	AcodePhoenix AnalyteCriteria\$8260GWRAcetoneMA / CAM Protocol / VOA AQ RL\$8260GWRCarbon DisulfideMA / CAM Protocol / VOA AQ RL\$8260GWRTetrahydrofuran (THF)MA / CAM Protocol / VOA AQ RL\$8260GWRtrans-1,4-dichloro-2-buteneMA / CAM Protocol / VOA AQ RL\$8260GWR1,2-DibromoethaneMA / CAM Protocol / GW-1 (mg/l)\$8260GWR1,2-DibromoethaneMA / GROUNDWATER STANDARDS / GW-1\$8270-SIMFSRBenzoic acidMA / CAM Protocol / SVOA AQ RLAS-WMArsenicMA / CAM R 310.40.1600 / GW-1 (mg/l)AS-WMArsenicMA / GROUNDWATER STANDARDS / GW-1\$8260GWRCarbon DisulfideMA / CAM Protocol / VOA AQ RL\$8260GWRTetrahydrofuran (THF)MA / CAM Protocol / VOA AQ RL\$8260GWRtrans-1,4-dichloro-2-buteneMA / CAM Protocol / VOA AQ RL\$8260GWR1,2-DibromoethaneMA / CAM Protocol / VOA AQ RL\$8260GWR1,2-DibromoethaneMA / CAM R 310.40.1600 / GW-1 (mg/l)\$8260GWR1,2-DibromoethaneMA / GROUNDWATER STANDARDS / GW-1\$MCPADD-WM1,4-DioxaneMA / CMR 310.40.1600 / GW-1 (mg/l)	Acode Phoenix Analyte Criteria Result \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND \$8260GWR trans-1,4-dichloro-2-butene MA / CAM Protocol / VOA AQ RL ND \$8260GWR 1,2-Dibromoethane MA / GROUNDWATER STANDARDS / GW-1 ND \$8270-SIMFSR Benzoic acid MA / CAM Protocol / SVOA AQ RL ND AS-WM Arsenic MA / CMR 310.40.1600 / GW-1 (mg/l) 0.011 AS-WM Arsenic MA / GROUNDWATER STANDARDS / GW-1 0.011 \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND	Acode Phoenix Analyte Criteria Result RL \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND 25 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND 2.5 \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR trans-1,4-dichloro-2-butene MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR 1,2-Dibromoethane MA / GROUNDWATER STANDARDS / GW-1 ND 0.50 \$8270-SIMFSR Benzoic acid MA / CAM Protocol / SVOA AQ RL ND 48 AS-WM Arsenic MA / CAM Protocol / GW-1 (mg/l) 0.011 0.004 AS-WM Arsenic MA / GROUNDWATER STANDARDS / GW-1 0.011 0.004 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR Tetrah	Acode Phoenix Analyte Criteria Result RL Criteria \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND 25 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND 2.5 \$8260GWR trans-1,4-dichloro-2-butene MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR 1,2-Dibromoethane MA / CAM Protocol / GW-1 (mg/l) ND 0.50 0.02 \$8260GWR 1,2-Dibromoethane MA / GROUNDWATER STANDARDS / GW-1 ND 0.50 0.02 \$8270-SIMFSR Benzoic acid MA / CAM Protocol / SVOA AQ RL ND 48 AS-WM Arsenic MA / CAM Protocol / GW-1 (mg/l) 0.011 0.004 0.01 \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND 25 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND<	Acode Phoenix Analyte Criteria Criteria REsult RL Criteria Criteria \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND 25 10 \$8260GWR Carbon Disulfide MA / CAM Protocol / VOA AQ RL ND 5.0 2 \$8260GWR Tetrahydrofuran (THF) MA / CAM Protocol / VOA AQ RL ND 5.0 2 \$8260GWR trans-1,4-dichloro-2-butene MA / CAM Protocol / VOA AQ RL ND 5.0 2 \$8260GWR 1,2-Dibromoethane MA / CAM Protocol / VOA AQ RL ND 0.50 0.02 0.02 \$8260GWR 1,2-Dibromoethane MA / GROUNDWATER STANDARDS / GW-1 ND 0.50 0.02 0.02 \$8270-SIMFSR Benzoic acid MA / CAM Protocol / SVOA AQ RL ND 48 10 AS-WM Arsenic MA / CAM Protocol / GW-1 (mg/l) 0.011 0.004 0.01 0.01 \$8260GWR Acetone MA / CAM Protocol / VOA AQ RL ND 25 10 \$8260GWR Carbon Disulfide

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.

MassDEP Analytical Protocol Certification Form										
Laboratory Name: Phoenix Environmental Laboratories, Inc. Project #:										
Project Location: 30 PENNIMAN RTN:										
This Form provides certifications for the following data set: [list Laboratory Sample ID Number(s)]										
CF10995, CF11044										
Matrices of Commitment Confees Water Co. 2010 attention Co. 2010 atten										
Matrices: ✓ Groundwater/Surface Water Soil/Sediment Drinking Water Air ✓ Other: WATER, W CAM Protocol (check all that apply below)										
8260 VOC 7470/7471 Hg MassDEP VPH 8081 Pesticides 7196 Hex Cr MassDEP APH										
CAM II A		CAM III B	V	CAM IV A	CAM V B		CAM VI B		CAM I	X A 🗌
8270 SVOC CAM II B		7010 Metals CAM III C		MassDEP EPH CAM IV B	8151 Herbicides CAM V C		8330 Explosives CAM VIII A		TO-15 CAM I	
6010 Metals CAM III A		6020 Metals CAM III D		8082 PCB CAM V A	9012 Total Cyanide/PAC CAM V1 A		6860 Perchlorat CAM VIII B	е		
Affirmative responses to questions A through F are required for "Presumptive Certainty" status										
Α	Were all samples received in a condition consistent with those described on the Chain-of-Custody, properly preserved (including temperature*) in the field or laboratory, and prepared/analyzed with method holding times? (* see narrative) ✓ Yes								□ No	
В	Were the analytical method(s) and all associated QC requirements specified in the selected CAM protocol(s) followed?							\	Yes	□No
С	Were all required corrective actions and analytical response actions specified in the selected CAM protocol(s) implemented for all identified performance standard non-conformances?								Yes	□No
D	Does the laboratory report comply with all the reporting requirements speified in CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data"?								Yes	□No
Е								□No		
b. APH and TO-15 methods only: Was the complete analyte method?					ete analyte list	repor	ted for each		Yes	□ No
F	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)? ✓ Yes □ N								□ No	
Responses to questions G, H and I below is required for "Presumptive Certainty" status										
G	Were the reporting limits at or below all CAM reporting limits specified in the selected CAM protocol(s)?								Yes	✓ No
Data User Note: Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40. 1056(2)(k) and WSC-07-350										
Н	See Secti	on: SVOASI	M Narra						Yes	☑ No
I	protocol(s)?								Yes	✓ No
All negative responses must be addressed in an attached laboratory narrative.										
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.										
Date: Thursday, January 23, 2020										
Authorized Roshui Waket Printed Name: Rashmi Maket						•	•			
Signature:				– Posit	Position: Project Manager					



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



MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

SDG Comments

Metals Analysis:

The client requested a site specific list of elements which is shorter than the 6010 MCP list.

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.

504.1

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

Instrument:

CHEM35 01/16/20-1

Chelsey Guerette, Chemist 01/16/20

CF10995 (1X)

The initial calibration (CHEM35/504tcp_0116): 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 15% except for the following compounds:None.

QC (Batch Specific):

Batch 514515 (CF13729)

CF10995

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 25% with the following exceptions: None.

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.

Cyanide Narration

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

Instrument:

LACHAT 01/16/20-2

Eric Geyer, Greg Danielewski, Chemist 01/16/20

CF10995

The samples were distilled in accordance with the method.

The initial calibration met criteria.

The calibration check standards (ICV,CCV) were within 15% of true value and were analyzed at a frequencey of one per ten samples.

The continuing calibration blanks (ICB,CCB) had concentrations less than the reporting level.

The method blank, laboratory control sample (LCS), and matrix spike were distilled with the samples.

QC (Batch Specific):



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

Cyanide Narration

Batch 514435 (CF10215)

CF10995

All LCS recoveries were within 90 - 110 with the following exceptions: None.

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

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.

Hexavalent Chromium (Aqueous)

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

Instrument:

BECKMAN DU720 01/13/20-1 Dustin Harrison, Chemist 01/13/20

CF10995

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

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

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

QC (Batch Specific):

Batch 513988 (CF11334)

CF10995

All LCS recoveries were within 90 - 110 with the following exceptions: None.

Additional Hexavalent Chromium criteria: LCS acceptance range for waters is 90-110% and MS acceptance range is 85-115%.

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 01/15/20 07:38 Rick Schweitzer, Chemist 01/15/20

CF10995

The method preparation blank contains all of the acids and reagents as the samples; the instrument blanks do not.

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):



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Certification Report

January 23, 2020 SDG I.D.: GCF10995

Mercury Narration

Batch 514110 (CF11603)

CF10995

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 01/14/20 08:53

Emily Kolominskaya, Chemist 01/14/20

CF10995

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 514163 (CF10467)

CF10995

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

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

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

LACHAT

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

Instrument:

LACHAT 01/13/20-1

Thomas Budz, Chemist 01/13/20

CF10995

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

All method verification standards and blanks met criteria.

QC (Batch Specific):

Batch 514023 (CF10330)

CF10995

All LCS recoveries were within 90 - 110 with the following exceptions: None.



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

LACHAT

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.

NITROGEN

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

Instrument:

LACHAT 01/16/20-1

Kandi Della Bella, Chemist 01/16/20

CF10995

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

All method verification standards and blanks met criteria.

QC (Batch Specific):

Batch 514457 (CF10393)

CF10995

All LCS recoveries were within 85 - 115 with the following exceptions: None.

Additional criteria: LCS acceptance range for waters is 85-115% and for soils is 75-125%. MS acceptance range is 75-125%.

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.

PCB Narration

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

Instrument:

AU-ECD24 01/15/20-1

Saadia Chudary, Chemist 01/15/20

CF10995 (1X)

The initial calibration (PC1125AI) RSD for the compound list was less than 20% except for the following compounds: None.

The initial calibration (PC1125BI) 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 15% except for the following compounds:None.

QC (Batch Specific):

Batch 514154 (CF10995)

CF10995

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

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

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

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

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.



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

SVOA Narration

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

Instrument:

CHEM06 01/15/20-1

Wes Bryon, Chemist 01/15/20

CF10995 (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 (CHEM06/6_SPLIT_0108):

100% of target compounds met criteria.

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

The following compounds did not meet recommended response factors: 2-Nitrophenol 0.076 (0.1)

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

Continuing Calibration Verification (CHEM06/0115_03-6_SPLIT_0108) (MCP Compliance):

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

95% 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 0.079 (0.1)

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

QC (Batch Specific):

Batch 513952 (CF10667)

CF10995

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

This batch consists of a Blank and LCS

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.

SVOA-Dioxane

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

Instrument:

CHEM22 01/15/20-1

Adam Werner, Chemist 01/15/20

CF10995 (1X)

Initial Calibration Evaluation (CHEM22/DIOX_0110):

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.



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

SVOA-Dioxane

Continuing Calibration Verification (CHEM22/0115_05-DIOX_0110) (MCP Compliance):

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: None.

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

QC (Batch Specific):

Batch 514189 (CF12265)

CF10995

All LCS recoveries were within 30 - 130 with the following exceptions: None.

All LCSD recoveries were within 30 - 130 with the following exceptions: None.

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

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

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

QC Batch 513952 (Samples: CF10995): -----

The blank surrogate was below criteria. (% Nitrobenzene-d5(CF10667))

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)

Instrument:

CHEM27 01/15/20-2

Wes Bryon, Chemist 01/15/20

CF10995 (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_0103):

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 (CHEM27/0115_11-27_SIM18_0103) (MCP Compliance):

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: None.

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



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

SVOASIM Narration

QC (Batch Specific):

Batch 513952 (CF10667)

CF10995

All LCS recoveries were within 40 - 140 with the following exceptions: Hexachlorocyclopentadiene(33%)

This batch consists of a Blank and LCS

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? Yes.

Instrument:

CHEM17 01/13/20-2

Michael Hahn, Chemist 01/13/20

CF11044 (1X)

Initial Calibration Evaluation (CHEM17/VT-S011220):

98% of target compounds met criteria.

The following compounds had %RSDs >20%: 1,2-Dibromo-3-chloropropane 25% (20%), trans-1,4-dichloro-2-butene 32% (20%) The following compounds did not meet recommended response factors: 1,2-Dibromo-3-chloropropane 0.035 (0.05), 2-Hexanone 0.087 (0.1), Acetone 0.052 (0.1), Bromoform 0.094 (0.1), Methyl ethyl ketone 0.078 (0.1), Tetrahydrofuran (THF) 0.046 (0.05), trans-1,4-dichloro-2-butene 0.041 (0.05)

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

Continuing Calibration Verification (CHEM17/0113_30-VT-S011220) (MCP Compliance):

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

95% 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: 1,2-Dibromo-3-chloropropane 0.032 (0.05), 2-Hexanone 0.077 (0.1), Acetone 0.058 (0.1), Bromoform 0.092 (0.1), Methyl ethyl ketone 0.070 (0.1), Tetrahydrofuran (THF) 0.043 (0.05), trans-1,4-dichloro-2-butene 0.031 (0.05)

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

CHEM17 01/14/20-2

Michael Hahn, Chemist 01/14/20

CF10995 (1X)

Initial Calibration Evaluation (CHEM17/VT-S011220):

98% of target compounds met criteria.

The following compounds had %RSDs >20%: 1,2-Dibromo-3-chloropropane 25% (20%), trans-1,4-dichloro-2-butene 32% (20%) The following compounds did not meet recommended response factors: 1,2-Dibromo-3-chloropropane 0.035 (0.05), 2-Hexanone 0.087 (0.1), Acetone 0.052 (0.1), Bromoform 0.094 (0.1), Methyl ethyl ketone 0.078 (0.1), Tetrahydrofuran (THF) 0.046 (0.05), trans-1,4-dichloro-2-butene 0.041 (0.05)

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

Continuing Calibration Verification (CHEM17/0114_31-VT-S011220) (MCP Compliance):



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

VOA Narration

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

95% 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: 1,2-Dibromo-3-chloropropane 0.033 (0.05), 2-Hexanone 0.075 (0.1), Acetone 0.041 (0.1), Bromoform 0.089 (0.1), Methyl ethyl ketone 0.063 (0.1), Tetrahydrofuran (THF) 0.040 (0.05), trans-1,4-dichloro-2-butene 0.043 (0.05)

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

QC (Batch Specific):

Batch 514113 (CF10770) CHEM17 1/13/2020-2

CF11044

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: Naphthalene(31.1%)

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

Batch 514336 (CF12269) CHEM17 1/14/2020-2

CF10995

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%.

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-OXY Narration

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

Instrument:

CHEM17 01/15/20-1 Michael Hahn, Chemist 01/15/20

CF10995 (1X)

Initial Calibration Evaluation (CHEM17/OXY010720):

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 (CHEM17/0115_05-OXY010720) (MCP Compliance):

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: None.



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MCP Certification Report

January 23, 2020 SDG I.D.: GCF10995

VOA-OXY Narration

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

QC (Batch Specific):

Batch 514382 (CF10995) CHEM17 1/15/2020-1

CF10995

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 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.

Stank * SURCHARGE APPLIES V Phoenix Std Report This section MUST be Cooler Yes No Full Data Package Bottle Quantities. ☐ Tier II Checklist completed with Data Delivery/Confact Options: 9560189 Data Package Richal, leany o GIS/Key Data Format EQuIS Other S-2 GW-1 S-2 GW-2 S-2 GW-3 S-1 GW-1 S-1 GW-2 S-1 GW-3 S-3 GW-1 S-3 GW-2 S-3 GW-3 Project P.O. Coolant: ☐ MWRA eSMART Phone: と Email: MA CP Certification SW Protection GW-1 GW-2 Residential DEC 587 East Middle Turnpike, P.O. Box 370, Manchester, CT 06040 ☐ GW Protection SW Protection GA Mobility **GB Mobility** CI RCP Cert ST (21819 BA Fax (860) 645-0823 I/C DEC 50 Persona Ro **CHAIN OF CUSTODY RECORD** Kidhar Leurad Client Services (860) 645-8726 (Comm/Industrial) GA Leachability GB Leachability Direct Exposure Direct Exposure GA-GW Objectives (Residential) Objectives GB-GW Email: info@phoenixlabs.com Report to: Invoice to: QUOTE# 교 Project: Analysis Request * SURCHARGE APPLIES 25.3 Turnaround Time: Time: Standard 3 Days* 2 Days* 1115120 ☐ 1 Day* 04/1/ Sampled Time 12:00 1110/20 32.50 Matrix Code:

DW=Drinking Water GW=Ground Water SW=Surface Water WW=Waste Water RW=Raw Water SE=Sediment SL=Sludge S=Soil SD=Soild W=Wipe OIL=Oil B=Bulk L=Liquid X = ____(Other) Date: Date Sampled 2 2 Client Sample - Information - Identification * OI Diss BORTHAL FIELL, hold will the Sample Matrix <u>ک</u> ટુ Environmental Laboratories, Inc. Comments, Special Requirements or Regulations: Accepted by: Customer Sample Identification Quiney MA $O_{i,s}$ Storte £ PHOENIX USE ONLY Relinquished by Customer: Address: 2660 SAMPLE # 1000 Sampler's Signature

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Phoenix Environmental Laboratories, Inc. 587 East Middle Turnpike
Manchester, CT 06040
(860) 812-0086
Please call Lisa with any questions

Container Order

Company: Stantec - Quincy, MA

RGP Project: Contact:

Ryan Willis 1/8/20 Date:

Ground Water

<u>Parameter</u>	Ammonia	Chloride, Chlorine, TSS, Hex Cr, Tri Cr	Sb, As, Cd, Cu, Fe, Pb, Hg, Ni, Se, Ag, Zn	Cyanide	VOC	1,4 Dioxane	SVOC	PCB	TPH 1164	Alcohols	Oxygenates	EDB
Preservative	H2SO4	AS IS	HN03	NAOH	HCL	NAHSO4	AS IS	AS IS	H2S04	AS IS	TSP	AS IS
Container	250ml Plastic	500ml Plastic	250ml Plastic	250ml Plastic	40ml Vials	8oz Amber	Liter Amber	Liter Amber	Liter Amber	40ml Vials	40ml Vials	40ml Vials
<u>Total</u>	-	τ-	τ	ν	က	_	2	-	-	2	2	2
Sets	_	-	_	_	_	_	_	_	_	_	_	_
# Per Set	-	_	_		က	_	7	_	τ-	7	7	7

Also Included:

CT Chain, Labels 2 HCL Vials with Reagent Water wrapped = Trip Blank

Makrina Nolan

Subject:

GCF10995

From: Learned, Richard [mailto:Richard.Learned@stantec.com]

Sent: Tuesday, January 14, 2020 11:10 AM

To: Makrina Nolan

Subject: Re: Samples received yesterday

Yes please run. This sample is labeled D1.

Richard Learned LSP

Senior Environmental Project Manager

Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stanted

146 Main Street Unit 3 Hyannis MA 02601-3128 US



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From: Makrina Nolan < Makrina@phoenixlabs.com>

Sent: Tuesday, January 14, 2020 11:05 AM

To: Learned, Richard < Richard.Learned@stantec.com>

Subject: Samples received yesterday

Good morning,

We received your samples yesterday, with regards to the attached chain. Unfortunately, the first sample "DI" was received past hold for HexChrome.

Please let me know if you would like to run HexChrome past hold for this sample.

Thank you,

Makrina Nolan Client Service –Project Manager Drinking Water Specialist Phoenix Environmental Labs 587 Middle Turnpike East Manchester, CT

Direct Line: 860-645-3219

Website: www.phoenixlabs.com



Friday, February 21, 2020

Attn: Mr Joseph Salvetti

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Project ID: 195601892 SDG ID: GCF32692

Sample ID#s: CF32700 - CF32703, CF32712 - CF32713

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

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

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

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

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

Sincerely yours,

Phyllis/Shiller

Laboratory Director

NELAC - #NY11301

CT Lab Registration #PH-0618
MA Lab Registration #M-CT007

ME Lab Registration #CT-007

NH Lab Registration #213693-A,B

NJ Lab Registration #CT-003 NY Lab Registration #11301

PA Lab Registration #68-03530 RI Lab Registration #63

UT Lab Registration #CT00007 VT Lab Registration #VT11301



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



SDG Comments

February 21, 2020

SDG I.D.: GCF32692

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



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Sample Id Cross Reference

February 21, 2020

SDG I.D.: GCF32692

Project ID: 195601892

Client Id	Lab Id	Matrix
B1	CF32700	SOIL
B2	CF32701	SOIL
B3	CF32702	SOIL
B4	CF32703	SOIL
C4	CF32712	WATER
D1	CF32713	WATER



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

Analysis Report

February 21, 2020

FOR: Attn: Mr Joseph Salvetti

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample InformationCustody InformationDateTimeMatrix:WATERCollected by:02/14/2013:30Location Code:STANTECMAReceived by:LB02/14/2017:16

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

P.O.#: Laboratory Data

SDG ID: GCF32692

Phoenix ID: CF32712

Project ID: 195601892

Client ID: C4

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Calcium	168	0.10	mg/L	10	02/18/20	CPP	SW6010D/E200.7
Hardness (CaCO3)	567	0.1	mg/L	1	02/18/20		E200.7
Magnesium	35.7	0.010	mg/L	1	02/18/20	MGH	SW6010D/E200.7
Total Metals Digestion	Completed				02/17/20	AG	

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

February 21, 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

February 21, 2020

FOR: Attn: Mr Joseph Salvetti

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample InformationCustody InformationDateTimeMatrix:WATERCollected by:02/14/2013:45Location Code:STANTECMAReceived by:LB02/14/2017:16

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

P.O.#: Laboratory Data

SDG ID: GCF32692

Phoenix ID: CF32713

Project ID: 195601892

Client ID: D1

Parameter	Result	RL/ PQL	Units	Dilution	Date/Time	Ву	Reference
Calcium	140	0.010	mg/L	1	02/18/20	MGH	SW6010D/E200.7
Hardness (CaCO3)	470	0.1	mg/L	1	02/18/20		E200.7
Magnesium	29.2	0.010	mg/L	1	02/18/20	MGH	SW6010D/E200.7
Total Metals Digestion	Completed				02/17/20	AG	

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

February 21, 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

February 21, 2020

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

QA/QC Data

SDG I.D.: GCF32692

J ,													
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 518723 (mg/kg)), QC San	nple No:	CF3217	8 2X (CF	32700	, CF327	701, CF3	32702,	CF3270	03)			
Mercury - Soil	BRL	0.03	< 0.03	< 0.03	NC	110	97.6	11.9	94.7	94.5	0.2	75 - 125	20
Comment:													
Additional Mercury criteria: LCS	acceptano	e range t	for waters	is 80-120	% and fo	or soils i	s 75-125°	%					
QA/QC Batch 519031 (mg/L),	QC Sam	ple No: (CF33257	(CF327	03)								
Mercury - Water	BRL	0.0002	< 0.0002	<0.0002	NC	111			117			75 - 125	30
Comment:													
Additional Mercury criteria: LCS	acceptano	e range t	for waters	is 80-120	% and fo	or soils i	s 75-125°	%					
QA/QC Batch 518805 (mg/L),	QC Sam	ple No: (CF31592	(CF327	12, CF3	32713)							
ICP Metals - Aqueous													
Calcium	BRL	0.010	0.153	0.149	2.60	105	105	0.0	105			80 - 120	20
Magnesium	BRL	0.010	0.043	0.042	NC	104	104	0.0	105			80 - 120	20
Comment:													



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QA/QC Report

February 21, 2020

QA/QC Data

SDG I.D.: GCF32692

<i>y</i> ,		-									
Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
QA/QC Batch 518767H (mg/K)	g), QC Sa	imple No: CF31429 50X (CF32	703 (50	X))							
Volatiles - Soil (High Lev	_	•	`	, ,							
Tetrachloroethene	<u>u.,</u> ND	0.25	117	119	1.7	118	112	5.2	70 - 130	30	
% 1,2-dichlorobenzene-d4	100	%	101	102	1.0	101	100	1.0	70 - 130	30	
% Bromofluorobenzene	96	%	99	100	1.0	100	101	1.0	70 - 130	30	
% Dibromofluoromethane	95	%	97	96	1.0	94	92	2.2	70 - 130	30	
% Toluene-d8	97	%	99	99	0.0	98	99	1.0	70 - 130	30	
Comment:											
Additional 8260 criteria: 10% of o	compounds	can be outside of acceptance crite	eria as lo	ng as rec	overy is	10%.					
	-	nple No: CF31907 (CF32700, C		_	_						
Volatiles - Soil (Low Leve		.,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	. 02700	• •							
1,1,1,2-Tetrachloroethane	ND	0.005	105	110	4.7	92	97	5.3	70 - 130	30	
1,1,1-Trichloroethane	ND	0.005	96	98	2.1	91	94	3.2	70 - 130	30	
1,1,2,2-Tetrachloroethane	ND	0.003	105	110	4.7	82	84	2.4	70 - 130	30	
1,1,2-Trichloroethane	ND	0.005	102	105	2.9	88	92	4.4	70 - 130	30	
1,1-Dichloroethane	ND	0.005	96	99	3.1	91	97	6.4	70 - 130	30	
1,1-Dichloroethene	ND	0.005	103	102	1.0	93	100	7.3	70 - 130	30	
1,1-Dichloropropene	ND	0.005	101	100	1.0	90	92	2.2	70 - 130	30	
1,2,3-Trichlorobenzene	ND	0.005	117	121	3.4	41	40	2.5	70 - 130	30	m
1,2,3-Trichloropropane	ND	0.005	105	111	5.6	84	86	2.4	70 - 130	30	
1,2,4-Trichlorobenzene	ND	0.005	116	117	0.9	45	44	2.2	70 - 130	30	m
1,2,4-Trimethylbenzene	ND	0.001	108	109	0.9	74	78	5.3	70 - 130	30	
1,2-Dibromo-3-chloropropane	ND	0.005	106	118	10.7	73	77	5.3	70 - 130	30	
1,2-Dibromoethane	ND	0.005	104	108	3.8	84	88	4.7	70 - 130	30	
1,2-Dichlorobenzene	ND	0.005	102	105	2.9	56	62	10.2	70 - 130	30	m
1,2-Dichloroethane	ND	0.005	99	102	3.0	90	93	3.3	70 - 130	30	
1,2-Dichloropropane	ND	0.005	100	102	2.0	91	94	3.2	70 - 130	30	
1,3,5-Trimethylbenzene	ND	0.001	107	109	1.9	82	85	3.6	70 - 130	30	
1,3-Dichlorobenzene	ND	0.005	105	106	0.9	66	67	1.5	70 - 130	30	m
1,3-Dichloropropane	ND	0.005	103	107	3.8	89	93	4.4	70 - 130	30	
1,4-Dichlorobenzene	ND	0.005	102	104	1.9	62	64	3.2	70 - 130	30	m
1,4-dioxane	ND	0.1	120	115	4.3	103	102	1.0	40 - 160	30	
2,2-Dichloropropane	ND	0.005	100	95	5.1	90	89	1.1	70 - 130	30	
2-Chlorotoluene	ND	0.005	103	106	2.9	69	77	11.0	70 - 130	30	m
2-Hexanone	ND	0.025	95	102	7.1	73	81	10.4	40 - 160	30	
2-Isopropyltoluene	ND	0.005	104	105	1.0	74	78	5.3	70 - 130	30	
4-Chlorotoluene	ND	0.005	102	103	1.0	69	73	5.6	70 - 130	30	m
4-Methyl-2-pentanone	ND	0.025	99	103	4.0	88	93	5.5	40 - 160	30	
Acetone	ND	0.01	72	81	11.8	57	65	13.1	40 - 160	30	
Acrylonitrile	ND	0.005	93	97	4.2	76	86	12.3	70 - 130	30	
Benzene	ND	0.001	105	105	0.0	95	97	2.1	70 - 130	30	
Bromobenzene	ND	0.005	105	109	3.7	75	77	2.6	70 - 130	30	

QA/QC Data

SDG I.D.: GCF32692

Parameter	Blank	Blk RL	LCS %	LCSD %	LCS RPD	MS %	MSD %	MS RPD	% Rec Limits	% RPD Limits	
Bromochloromethane	ND	0.005	99	104	4.9	89	90	1.1	70 - 130	30	
Bromodichloromethane	ND	0.005	105	107	1.9	92	95	3.2	70 - 130	30	
Bromoform	ND	0.005	108	116	7.1	78	86	9.8	70 - 130	30	
Bromomethane	ND	0.005	112	111	0.9	108	110	1.8	40 - 160	30	
Carbon Disulfide	ND	0.005	103	105	1.9	88	93	5.5	70 - 130	30	
Carbon tetrachloride	ND	0.005	101	102	1.0	89	95	6.5	70 - 130	30	
Chlorobenzene	ND	0.005	104	107	2.8	82	86	4.8	70 - 130	30	
Chloroethane	ND	0.005	92	93	1.1	89	87	2.3	70 - 130	30	
Chloroform	ND	0.005	96	100	4.1	89	95	6.5	70 - 130	30	
Chloromethane	ND	0.005	83	84	1.2	76	80	5.1	40 - 160	30	
cis-1,2-Dichloroethene	ND	0.005	104	95	9.0	87	84	3.5	70 - 130	30	
cis-1,3-Dichloropropene	ND	0.005	104	107	2.8	83	88	5.8	70 - 130	30	
Dibromochloromethane	ND	0.003	113	119	5.2	91	98	7.4	70 - 130	30	
Dibromomethane	ND	0.005	101	105	3.9	84	88	4.7	70 - 130	30	
Dichlorodifluoromethane	ND	0.005	83	83	0.0	78	83	6.2	40 - 160	30	
Diethyl ether	ND	0.005	86	92	6.7	104	108	3.8	70 - 130	30	
Di-isopropyl ether	ND	0.005	88	93	5.5	87	93	6.7	70 - 130	30	
Ethyl tert-butyl ether	ND	0.005	92	98	6.3	91	97	6.4	70 - 130	30	
Ethylbenzene	ND	0.001	108	109	0.9	88	93	5.5	70 - 130	30	
Hexachlorobutadiene	ND	0.005	114	110	3.6	55	58	5.3	70 - 130	30	m
Isopropylbenzene	ND	0.001	106	108	1.9	86	89	3.4	70 - 130	30	
m&p-Xylene	ND	0.002	108	108	0.0	84	91	8.0	70 - 130	30	
Methyl ethyl ketone	ND	0.005	82	86	4.8	72	79	9.3	40 - 160	30	
Methyl t-butyl ether (MTBE)	ND	0.001	94	100	6.2	92	99	7.3	70 - 130	30	
Methylene chloride	ND	0.005	87	92	5.6	84	89	5.8	70 - 130	30	
Naphthalene	ND	0.005	126	132	4.7	<10	34	NC	70 - 130	30	l,m
n-Butylbenzene	ND	0.001	114	110	3.6	70	74	5.6	70 - 130	30	.,
n-Propylbenzene	ND	0.001	108	106	1.9	79	83	4.9	70 - 130	30	
o-Xylene	ND	0.002	107	109	1.9	82	89	8.2	70 - 130	30	
p-Isopropyltoluene	ND	0.001	112	111	0.9	78	81	3.8	70 - 130	30	
sec-Butylbenzene	ND	0.001	114	114	0.0	79	84	6.1	70 - 130	30	
Styrene	ND	0.005	108	111	2.7	62	75	19.0	70 - 130	30	m
tert-amyl methyl ether	ND	0.005	97	102	5.0	95	98	3.1	70 - 130	30	
tert-Butylbenzene	ND	0.001	107	107	0.0	81	84	3.6	70 - 130	30	
Tetrachloroethene	ND	0.005	106	105	0.9	91	91	0.0	70 - 130	30	
Tetrahydrofuran (THF)	ND	0.005	88	95	7.7	79	87	9.6	70 - 130	30	
Toluene	ND	0.001	106	106	0.0	90	94	4.3	70 - 130	30	
trans-1,2-Dichloroethene	ND	0.005	102	104	1.9	90	96	6.5	70 - 130	30	
trans-1,3-Dichloropropene	ND	0.005	105	107	1.9	77	81	5.1	70 - 130	30	
trans-1,4-dichloro-2-butene	ND	0.005	106	111	4.6	71	74	4.1	70 - 130	30	
Trichloroethene	ND	0.005	102	101	1.0	92	95	3.2	70 - 130	30	
Trichlorofluoromethane	ND	0.005	96	96	0.0	92	97	5.3	70 - 130	30	
Trichlorotrifluoroethane	ND	0.005	106	105	0.9	97	103	6.0	70 - 130	30	
Vinyl chloride	ND	0.005	94	95	1.1	89	93	4.4	70 - 130	30	
% 1,2-dichlorobenzene-d4	101	%	100	101	1.0	99	100	1.0	70 - 130	30	
% Bromofluorobenzene	97	%	101	102	1.0	99	101	2.0	70 - 130	30	
% Dibromofluoromethane	96	%	97	96	1.0	94	98	4.2	70 - 130	30	
% Toluene-d8	99	%	100	99	1.0	99	99	0.0	70 - 130	30	
Comment:	,,	.•	100	,,	1.5	,,	,,	0.0	, 5 - 130	30	

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

QA/QC Data

SDG I.D.: GCF32692

% % Blk LCS LCSD LCS MS **MSD** MS Rec **RPD** Blank RL **RPD** % RPD Limits % Limits Parameter QA/QC Batch 518756H (mg/Kg), QC Sample No: CF31907 50X (CF32701 (50X), CF32702 (50X)) Volatiles - Soil (High Level) 0.25 3.5 1,1,1,2-Tetrachloroethane 115 118 2.6 113 117 70 - 130 30 ND 1,1,1-Trichloroethane ND 0.25 108 105 2.8 108 107 0.9 70 - 130 30 ND 0.25 121 4.2 119 120 70 - 130 30 1,1,2,2-Tetrachloroethane 116 0.8 1,1,2-Trichloroethane ND 0.25 110 112 1.8 108 111 2.7 70 - 130 30 ND 107 70 - 130 1,1-Dichloroethane 0.25 107 0.0 110 108 1.8 30 70 - 130 1,1-Dichloroethene ND 0.25 103 103 0.0 103 101 2.0 30 ND 2.7 70 - 130 1,1-Dichloropropene 0.25 114 111 112 113 0.9 30 1,2,3-Trichlorobenzene ND 0.25 136 135 0.7 157 149 5.2 70 - 130 30 I,m 1,2,3-Trichloropropane ND 0.25 118 122 3.3 113 114 0.9 70 - 130 30 1,2,4-Trichlorobenzene ND 0.25 140 139 0.7 154 145 70 - 130 30 6.0 I,m 1,2,4-Trimethylbenzene ND 0.25 121 119 1.7 123 125 1.6 70 - 130 30 ND 125 129 1,2-Dibromo-3-chloropropane 0.25 3.1 118 121 2.5 70 - 130 30 1,2-Dibromoethane ND 0.25 115 117 1.7 115 116 0.9 70 - 130 30 1,2-Dichlorobenzene ND 0.25 119 118 8.0 119 120 0.8 70 - 130 30 ND 106 107 110 110 1,2-Dichloroethane 0.25 0.9 0.0 70 - 130 30 ND 0.9 1,2-Dichloropropane 0.25 108 108 0.0 109 110 70 - 130 30 1,3,5-Trimethylbenzene ND 0.25 121 120 0.8 123 125 1.6 70 - 130 30 ND 1,3-Dichlorobenzene 0.25 121 120 8.0 120 123 2.5 70 - 130 30 1,3-Dichloropropane ND 0.25 115 116 0.9 116 116 0.0 70 - 130 30 1,4-Dichlorobenzene ND 0.25 121 118 2.5 120 120 0.0 70 - 130 30 ND 5 128 125 2.4 122 129 5.6 40 - 160 1,4-dioxane 30 2,2-Dichloropropane ND 0.25 110 103 6.6 100 101 1.0 70 - 130 30 ND 0.25 118 0.9 121 2-Chlorotoluene 117 118 2.5 70 - 130 30 2-Hexanone ND 1.3 105 109 3.7 110 107 2.8 40 - 160 30 ND 0.25 2-Isopropyltoluene 118 117 0.9 118 119 70 - 130 0.8 30 4-Chlorotoluene ND 0.25 116 114 1.7 118 120 1.7 70 - 130 30 ND 4-Methyl-2-pentanone 1.3 106 107 0.9 110 109 0.9 40 - 160 30 Acetone ND 0.5 71 73 2.8 75 75 0.0 40 - 160 30 Acrylonitrile ND 0.25 101 107 5.8 110 101 8.5 70 - 130 30 Benzene ND 0.25 117 116 0.9 116 118 1.7 70 - 130 30 Bromobenzene ND 0.25 117 116 0.9 116 120 3.4 70 - 130 30 ND 0.25 107 99 110 Bromochloromethane 110 2.8 10.5 70 - 130 30 Bromodichloromethane ND 0.25 110 109 0.9 108 110 1.8 70 - 130 30 Bromoform ND 0.25 113 117 3.5 102 107 4.8 70 - 130 30 Bromomethane ND 0.25 89 93 89 94 4.4 5.5 40 - 160 30 Carbon Disulfide ND 0.25 106 105 0.9 103 103 0.0 70 - 130 30 ND 109 Carbon tetrachloride 0.25 107 1.9 103 106 2.9 70 - 130 30 Chlorobenzene ND 0.25 120 119 0.8 118 120 1.7 70 - 130 30 Chloroethane ND 0.25 31 31 0.0 35 34 2.9 70 - 130 30 I,m ND 107 Chloroform 0.25 105 105 0.0 107 0.0 70 - 130 30 Chloromethane ND 0.25 95 92 3.2 94 92 2.2 40 - 160 30 95 97 cis-1,2-Dichloroethene ND 0.25 96 1.0 96 1.0 70 - 130 30 ND 0.25 112 0.0 70 - 130 cis-1,3-Dichloropropene 112 110 111 0.9 30 Dibromochloromethane ND 0.15 120 122 1.7 113 119 5.2 70 - 130 30 Dibromomethane ND 0.25 107 108 0.9 109 111 1.8 70 - 130 30 Dichlorodifluoromethane ND 0.25 87 88 90 91 1.1 1.1 40 - 160 30 Diethyl ether ND 57 57 0.25 0.0 61 60 70 - 130 1.7 30 I.m 95 98 Di-isopropyl ether ND 0.25 3.1 104 101 2.9 70 - 130 30 Ethyl tert-butyl ether ND 0.25 98 100 2.0 103 101 2.0 70 - 130 30 Ethylbenzene ND 0.25 125 124 8.0 123 125 1 6 70 - 130 30

QA/QC Data

% % Blk LCS **LCSD** LCS MS **MSD RPD** MS Rec Blank RL % **RPD** % % RPD Limits Limits % Parameter Hexachlorobutadiene ND 0.25 141 137 2.9 137 139 1.4 70 - 130 30 l,m Isopropylbenzene ND 0.25 121 119 1.7 119 123 3.3 70 - 130 30 m&p-Xylene ND 0.25 124 124 0.0 123 125 1.6 70 - 130 30 96 ND 0.25 97 99 2.0 97 1.0 40 - 160 30 Methyl ethyl ketone Methyl t-butyl ether (MTBE) ND 0.25 102 104 1.9 107 105 1.9 70 - 130 30 Methylene chloride ND 0.25 95 97 2.1 99 97 2.0 70 - 130 30 Naphthalene ND 0.25 146 147 0.7 175 161 8.3 70 - 130 30 I.m ND 133 1.5 133 134 n-Butylbenzene 0.25 135 0.7 70 - 130 30 I.m n-Propylbenzene ND 8.0 123 70 - 130 30 0.25 121 120 121 1.6 o-Xylene ND 0.25 122 123 0.8 121 124 2.4 70 - 130 30 p-Isopropyltoluene ND 0.25 128 127 8.0 127 130 2.3 70 - 130 30 ND 0.25 129 130 2.3 70 - 130 sec-Butylbenzene 131 1.5 133 30 l,m ND 0.25 123 123 0.0 123 123 0.0 70 - 130 30 Styrene ND tert-amyl methyl ether 0.25 105 106 0.9 106 107 0.9 70 - 130 30 tert-Butylbenzene 119 ND 0.25 119 0.0 120 122 1.7 70 - 130 30 Tetrachloroethene ND 0.25 122 120 1.7 120 119 8.0 70 - 130 30 ND Tetrahydrofuran (THF) 0.25 98 100 2.0 104 100 3.9 70 - 130 30 ND 0.25 119 117 1.7 117 119 Toluene 1.7 70 - 130 30 ND 116 0.9 119 1.7 70 - 130 trans-1,2-Dichloroethene 0.25 115 117 30 trans-1,3-Dichloropropene ND 0.25 112 111 0.9 109 109 0.0 70 - 130 30 trans-1,4-dichloro-2-butene ND 0.25 120 2.5 114 2.7 117 111 70 - 130 30 Trichloroethene ND 0.25 116 115 0.9 113 114 0.9 70 - 130 30 ND 0.25 27 27 29 29 Trichlorofluoromethane 0.0 0.0 70 - 130 30 I,m Trichlorotrifluoroethane ND 0.25 107 107 0.0 106 104 1.9 70 - 130 30 Vinyl chloride ND 0.25 104 105 1.0 105 106 0.9 70 - 130 30 101 99 99 % 1,2-dichlorobenzene-d4 % 100 100 0.0 0.0 70 - 130 30 % Bromofluorobenzene 96 % 101 101 0.0 101 101 0.0 70 - 130 30 % Dibromofluoromethane 92 % 93 94 96 2.1 94 1.1 70 - 130 30 % Toluene-d8 97 % 99 98 1.0 98 98 0.0 70 - 130 30 Comment:

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

SDG I.D.: GCF32692

February 21, 2020

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

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

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

Friday, February 21, 2020

Criteria: MA: CAM, S1

Sample Criteria Exceedances Report

GCF32692 - STANTECMA

State: N	ИΑ
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MA						RL	Analysis
Acode	Phoenix Analyte	Criteria	Result	RL	Criteria	Criteria	Units
\$8260MAR	trans-1,3-Dichloropropene	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	100	10	10	ug/Kg
\$8260MAR	Dibromochloromethane	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	100	5	5	ug/Kg
\$8260MAR	cis-1,3-Dichloropropene	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	100	10	10	ug/Kg
\$8260MAR	1,1,2,2-Tetrachloroethane	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	100	5	5	ug/Kg
\$8260MAR	Dibromochloromethane	MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1	ND	100	5	5	ug/Kg
\$8260MAR	1,1,2,2-Tetrachloroethane	MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1	ND	100	5	5	ug/Kg
\$8260MAR	trans-1,3-Dichloropropene	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	92	10	10	ug/Kg
\$8260MAR	Dibromochloromethane	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	92	5	5	ug/Kg
\$8260MAR	cis-1,3-Dichloropropene	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	92	10	10	ug/Kg
\$8260MAR	1,1,2,2-Tetrachloroethane	MA / CMR 310.40.1600 / S1 (mg/kg)	ND	92	5	5	ug/Kg
\$8260MAR	Dibromochloromethane	MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1	ND	92	5	5	ug/Kg
\$8260MAR	1,1,2,2-Tetrachloroethane	MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1	ND	92	5	5	ug/Kg
\$8260MAR	Tetrachloroethene	MA / CMR 310.40.1600 / S1 (mg/kg)	2500	280	1000	1000	ug/Kg
\$8260MAR	Tetrachloroethene	MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1	2500	280	1000	1000	ug/Kg
	\$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR \$8260MAR	\$8260MAR trans-1,3-Dichloropropene \$8260MAR Dibromochloromethane \$8260MAR 1,1,2,2-Tetrachloroethane \$8260MAR Dibromochloromethane \$8260MAR Dibromochloromethane \$8260MAR Dibromochloromethane \$8260MAR 1,1,2,2-Tetrachloroethane \$8260MAR trans-1,3-Dichloropropene \$8260MAR Dibromochloromethane \$8260MAR cis-1,3-Dichloropropene \$8260MAR cis-1,3-Dichloropropene \$8260MAR Dibromochloromethane \$8260MAR Dibromochloromethane \$8260MAR Dibromochloromethane \$8260MAR Tetrachloroethane	Acode Phoenix Analyte Criteria \$8260MAR trans-1,3-Dichloropropene MA / CMR 310.40.1600 / S1 (mg/kg) \$8260MAR Dibromochloromethane MA / CMR 310.40.1600 / S1 (mg/kg) \$8260MAR cis-1,3-Dichloropropene MA / CMR 310.40.1600 / S1 (mg/kg) \$8260MAR 1,1,2,2-Tetrachloroethane MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1 \$8260MAR Dibromochloromethane MA / SOIL S-1 STANDARDS / S-1 Soil & GW-1 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SOIL S-1 STANDARDS / S-1 Soil & GW-1 ND 100 5 5 \$8260MAR Dibromochloromethane MA / CMR 310.40.1600 / S1 (mg/kg) ND 92 10 10 \$8260MAR Dibromochloromethane MA / CMR 310.40.1600 / S1 (mg/kg) ND 92 5 5 \$8260MAR 1,1,2,2-Tetrachloroethane M

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.

		Ма	ssDE	P Analytical P	rotocol Cei	rtific	ation Forn	า		
Labo	ratory Na	ame: Phoe	nix En	vironmental Labora	atories, Inc. I	Proje	ct #:			
Proje	ct Locati	i on: 1956	01892		ı	RTN:				
This F	orm provid	les certification	ons for	the following data set	: [list Laborator	y San	ple ID Number(s)]		
CF327	'00, CF3270	01, CF32702, (CF3270	3						
Matric	es: Gro	undwater/Sur	face Wa	ater Soil/Sedimen	t Drinking	Wate	r 🗌 Air	✓ 0	ther:	WATER
CAM	Protocol (check all th	at app	ly below)						
8260 V CAM II		7470/7471 H CAM III B	g ••	MassDEP VPH CAM IV A	8081 Pesticides CAM V B		7196 Hex Cr CAM VI B		MassD CAM IX	DEP APH X A
8270 S CAM II		7010 Metals CAM III C		MassDEP EPH CAM IV B	8151 Herbicides CAM V C		8330 Explosives CAM VIII A		TO-15 CAM I	
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	е		
	Affirmati	ve respons	es to q	uestions A through	F are required	d for	Presumptive	Certai	inty" s	status
Α	Chain-of-0	Custody, pro	perly p	a condition consiste reserved (including to alyzed with method ho	emperature*) in	the f	eld or	V	Yes	□ No
В		analytical m CAM protoco		s) and all associated owed?	QC requiremen	its sp	ecified in the	V	Yes	□ No
С		CAM protoco		actions and analytica plemented for all ider				✓ ,	Yes	□ No
D	CAM VII A		ssuran	omply with all the repo ce and Quality Contro Pata"?				✓ ,	Yes	□ No
E	significant modificati	t modification ons).	n(s)? (ods only: Was each refer to the individual	method(s) for	a list o	of significant		Yes	□ No
	method?			only: Was the compl			ted for each		Yes	□ No
F	conformal		ed and	tocol QC and perforn evaluated in a labora rough E)?			ling all "No"	✓ ,	Yes	□ No
	Resp	onses to q	uestio	ns G, H and I below	is required for	r "Pre	sumptive Cer	tainty'	' statu	IS
G		reporting lim CAM protoco		r below all CAM repo	orting limits spe	cified	in the		Yes	☑ No
				resumptive Certainty" cribed in 310 CMR 40.				data us	sability	and
Н	See Secti	on: VOA Na	ration						Yes	✓ No
ı	Were resu protocol(s)?		complete analyte lis	· 				Yes	✓ No
respor	sible for o	d, attest unde	r the pa	esponses must be addi ains and penalties of p tion, the material conf	perjury that, bas	ed up	on my personal	inqui		
					Da	ate: F	Friday, Februa	ry 21	, 2020)
	orized	Ra	سماد	i nakol	Printed Na	me: F	Rashmi Makol			
Sign	ature: -	- 3			– Posit	ion: F	Project Manag	jer		



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



MCP Certification Report

February 21, 2020 SDG I.D.: GCF32692

SDG Comments

Metals Analysis:

The client requested a site specific list of elements which is shorter than the 6010 MCP list.

Volatile Comment: (CF32701, CF32702)

The client provided low level could not be analyzed due to physical interference from too much soil. The high level preserved vial requires a dilution resulting in elevated reporting levels.

Mercury Narration

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

Instrument:

MERLIN 02/17/20 09:49 Rick Schw

Rick Schweitzer, Chemist 02/17/20

CF32700, CF32701, CF32702, CF32703

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.

MERLIN 02/19/20 14:12

Rick Schweitzer, Chemist 02/19/20

CF32703

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 518723 (CF32178)

CF32700, CF32701, CF32702, CF32703

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

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

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

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

Batch 519031 (CF33257)

CF32703

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



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Certification Report

February 21, 2020 SDG I.D.: GCF32692

ICP Metals Narration

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

Instrument:

BLUE 02/17/20 13:55

Cindy Pearce, Mike Hornak, Chemist 02/17/20

CF32712. CF32713

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.

BLUE 02/18/20 08:31

Cindy Pearce, Mike Hornak, Chemist 02/18/20

CF32712

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 518805 (CF31592)

CF32712, CF32713

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%.

VOA Narration



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MCP Certification Report

February 21, 2020 SDG I.D.: GCF32692

VOA Narration

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

QC Batch 518756 (Samples: CF32700, CF32703): -----

The QC recoveries for one or more analytes is below the method criteria. A slight low bias is likely. (Naphthalene)

QC Batch 518756H: ----

The QC recoveries for one or more analytes is below the method criteria. A slight low bias is likely. (Chloroethane, Diethyl ether, Trichlorofluoromethane)

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. (1,2,3-Trichlorobenzene, 1,2,4-Trichlorobenzene, Hexachlorobutadiene, Naphthalene, n-Butylbenzene, sec-Butylbenzene)

Instrument:

CHEM31 02/14/20-2

Jane Li, Chemist 02/14/20

CF32700 (1X), CF32701 (50X), CF32702 (50X), CF32703 (1X)

Initial Calibration Evaluation (CHEM31/VT-L021220P):

94% of target compounds met criteria.

The following compounds had %RSDs >20%: 1,2-Dibromo-3-chloropropane 21% (20%), Acetone 32% (20%), Bromoform 22% (20%), Chloroethane 21% (20%), Naphthalene 27% (20%)

The following compounds did not meet Table 4 recommended minimum response factors: Acetone 0.094 (0.1), Bromoform 0.097 (0.1), Tetrachloroethene 0.177 (0.2)

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

Continuing Calibration Verification (CHEM31/0214_34-VT-L021220P) (MCP Compliance):

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

86% 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: Acetone 0.064 (0.1), Bromoform 0.091 (0.1), Tetrachloroethene 0.155 (0.2), Trichloroethene 0.197 (0.2)

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

CHEM31 02/16/20-1

Jane Li, Chemist 02/16/20

CF32703 (50X)

Initial Calibration Evaluation (CHEM31/VT-L021220P):

94% of target compounds met criteria.

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

The following compounds did not meet Table 4 recommended minimum response factors: Tetrachloroethene 0.177 (0.2)

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

Continuing Calibration Verification (CHEM31/0216_03-VT-L021220P) (MCP Compliance):

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 Table 4 recommended minimum response factors: Tetrachloroethene 0.184 (0.2)



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MCP Certification Report

February 21, 2020 SDG I.D.: GCF32692

VOA Narration

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

QC (Batch Specific):

Batch 518756 (CF31907) CHEM31 2/14/2020-2

CF32700(1X), CF32703(1X)

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

All LCSD recoveries were within 70 - 130 with the following exceptions: Naphthalene(132%)

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 518756H (CF31907) CHEM31 2/14/2020-2

CF32701(50X), CF32702(50X)

All LCS recoveries were within 70 - 130 with the following exceptions: 1,2,3-Trichlorobenzene(136%), 1,2,4-

Trichlorobenzene(140%), Chloroethane(31%), Diethyl ether(57%), Hexachlorobutadiene(141%), Naphthalene(146%), n-Butylbenzene(135%), sec-Butylbenzene(131%), Trichlorofluoromethane(27%)

All LCSD recoveries were within 70 - 130 with the following exceptions: 1,2,3-Trichlorobenzene(135%), 1,2,4-

Trichlorobenzene(139%), Chloroethane(31%), Diethyl ether(57%), Hexachlorobutadiene(137%), Naphthalene(147%), n-

Butylbenzene(133%), Trichlorofluoromethane(27%)

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 518767H (CF31429) CHEM31 2/16/2020-1

CF32703(50X)

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.

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

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.

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Environmental Laboratories, Inc.	oratories, Inc	11.5	Email:	Email: info@phoenixlabs.com Client Services (86	6	Fax (860) 645-0823 645-8726	Phone:		JOSCOP. Salvoh" a stanfor	
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			* SURCHAR	* SURCHARGE APPLIES	Objectives				* SURCHARGE APPLIES	_

Bobbi Aloisa

From: Learned, Richard kichard.learned@stantec.com

Sent: Monday, February 17, 2020 3:31 PM

To: Salvetti, Joseph; Bobbi Aloisa

Cc:Hanna, RittaSubject:Re: Rush -Allston

Also, run the hardness for the GW samples C4 and D1.

Richard Learned LSP

Senior Environmental Project Manager

Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stantec

146 Main Street Unit 3 Hyannis MA 02601-3128 US



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From: Learned, Richard < Richard.Learned@stantec.com >

Sent: Monday, February 17, 2020 3:28 PM

To: Salvetti, Joseph < <u>Joseph.Salvetti@stantec.com</u>>; Bobbi Aloisa < <u>bobbi@phoenixlabs.com</u>>

Cc: Hanna, Ritta < Ritta. Hanna@stantec.com>

Subject: Re: Rush -Allston

Yes just mercury for soil samples B1-B4.

Richard Learned LSP

Senior Environmental Project Manager

Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stanted

146 Main Street Unit 3 Hyannis MA 02601-3128 US



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From: Salvetti, Joseph < <u>Joseph.Salvetti@stantec.com</u>>

Sent: Monday, February 17, 2020 3:25 PM To: Bobbi Aloisa < bobbi@phoenixlabs.com >

Cc: Hanna, Ritta <Ritta.Hanna@stantec.com>; Learned, Richard <Richard.Learned@stantec.com>

Subject: RE: Rush -Allston

So sorry

I saw that just now Thank you

Joseph Salvetti LSP

Senior Associate

Direct: 508 591-4327 Mobile: 508 509-7393 Fax: 617 786-7962

Joseph.Salvetti@stantec.com

Stanted

400 Crown Colony Drive Suite 200

Quincy MA 02169-0982



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From: Bobbi Aloisa < bobbi@phoenixlabs.com> Sent: Monday, February 17, 2020 3:22 PM

To: Salvetti, Joseph < <u>Joseph.Salvetti@stantec.com</u>>

Cc: Hanna, Ritta < Ritta. Hanna@stantec.com >; Learned, Richard < Richard. Learned@stantec.com >

Subject: RE: Rush -Allston

Hg and VOCs are both checked off on the chain that was sent with results.

From: Salvetti, Joseph [mailto:Joseph.Salvetti@stantec.com]

Sent: Monday, February 17, 2020 3:21 PM

To: Bobbi Aloisa

Cc: Hanna, Ritta; Learned, Richard

Subject: RE: Rush -Allston

Hi Bobbi

All we wanted was Hg? Is that correct Rich?

Joe

Joseph Salvetti LSP

Senior Associate

Direct: 508 591-4327 Mobile: 508 509-7393 Fax: 617 786-7962

Joseph.Salvetti@stantec.com

Stantec

400 Crown Colony Drive Suite 200 Quincy MA 02169-0982



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From: Bobbi Aloisa < bobbi@phoenixlabs.com > Sent: Monday, February 17, 2020 12:56 PM

To: Salvetti, Joseph < Joseph. Salvetti@stantec.com >

Subject: Rush

Bobbi Aloisa

Vice President | Director of Client Services
Phoenix Environmental Laboratories, Inc.
587 East Middle Turnpike | Manchester, CT 06040
Direct Line: (860)-645-8728
www.phoenixlabs.com



Bobbi Aloisa

From: Bobbi Aloisa

Sent: Tuesday, February 18, 2020 1:43 PM

To: Learned, Richard
Cc: Bobbi Aloisa
Subject: RE: Rush

Will do, on 32703 Phoenix ID

From: Learned, Richard [mailto:Richard.Learned@stantec.com]

Sent: Tuesday, February 18, 2020 9:16 AM

To: Bobbi Aloisa Subject: Re: Rush

Bobbi,

As we discussed, please run TCLP for mercury on the B4 sample. One day TAT.

Richard Learned LSP

Senior Environmental Project Manager Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stantec 146 Main Street Unit 3 Hyannis MA 02601-3128 US

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From: Bobbi Aloisa <<u>bobbi@phoenixlabs.com</u>> Sent: Monday, February 17, 2020 4:11 PM

To: Learned, Richard < Richard. Learned@stantec.com >

Subject: FW: Rush

From: Bobbi Aloisa

Sent: Monday, February 17, 2020 12:56 PM
To: Salvetti, Joseph (<u>Joseph.Salvetti@stantec.com</u>)

Subject: Rush

Bobbi Aloisa

Vice President | Director of Client Services Phoenix Environmental Laboratories, Inc. 587 East Middle Turnpike | Manchester, CT 06040

Direct Line: (860)-645-8728 www.phoenixlabs.com





Tuesday, February 18, 2020

Attn: Mr Joseph Salvetti Stantec 400 Crown Colony Drive Suite 200

Quincy MA 02169

Project ID: CHARLES RIVER

SDG ID: GCF32691 Sample ID#s: CF32691

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,

Phyllis/Shiller

Laboratory Director

NELAC - #NY11301 CT Lab Registration #PH-0618

MA Lab Registration #M-CT007
ME Lab Registration #CT-007

NH Lab Registration #213693-A,B

NJ Lab Registration #CT-003 NY Lab Registration #11301 PA Lab Registration #68-03530

RI Lab Registration #63

UT Lab Registration #CT00007 VT Lab Registration #VT11301



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



SDG Comments

February 18, 2020

SDG I.D.: GCF32691

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



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Sample Id Cross Reference

February 18, 2020

SDG I.D.: GCF32691

Project ID: CHARLES RIVER

Client Id	Lab Id	Matrix
CHARLES SW	CF32691	SURFACE WATER



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

Analysis Report

February 18, 2020

FOR: Attn: Mr Joseph Salvetti

Stantec

400 Crown Colony Drive

Suite 200

Quincy MA 02169

Sample InformationCustody InformationDateTimeMatrix:SURFACE WATERCollected by:02/14/2014:20Location Code:STANTECMAReceived by:LB02/14/2017:16

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

Laboratory Data

SDG ID: GCF32691

Phoenix ID: CF32691

Project ID: CHARLES RIVER Client ID: CHARLES SW

P.O.#:

		RL/					
Parameter	Result	PQL	Units	Dilution	Date/Time	Ву	Reference
Silver	< 0.001	0.001	mg/L	1	02/15/20	CPP	SW6010D
Arsenic	< 0.004	0.004	mg/L	1	02/15/20	CPP	SW6010D
Cadmium	< 0.001	0.001	mg/L	1	02/15/20	CPP	SW6010D
Chromium	< 0.001	0.001	mg/L	1	02/15/20	CPP	SW6010D
Copper	< 0.005	0.005	mg/L	1	02/15/20	CPP	SW6010D
Iron	0.494	0.010	mg/L	1	02/15/20	CPP	E200.7
Hardness (CaCO3)	74.4	0.1	mg/L	1	02/15/20		E200.7
Mercury	< 0.0002	0.0002	mg/L	1	02/17/20	RS	SW7470A
Nickel	< 0.001	0.001	mg/L	1	02/15/20	CPP	SW6010D
Lead	< 0.002	0.002	mg/L	1	02/15/20	CPP	SW6010D
Antimony	< 0.005	0.005	mg/L	1	02/15/20	CPP	SW6010D
Selenium	< 0.010	0.010	mg/L	1	02/15/20	CPP	SW6010D
Trivalent Chromium	< 0.01	0.01	mg/L	1	02/17/20	BA	Calculation
Zinc	0.009	0.004	mg/L	1	02/15/20	CPP	SW6010D
Chromium, Hexavalent	< 0.01	0.01	mg/L	1	02/14/20 18:24	0	SM3500CRB-11
Ammonia as Nitrogen	0.14	0.05	mg/L	1	02/15/20	KDB	E350.1
рН	7.31	1.00	pH Units	1	02/14/20 22:08	AP	SM4500-H B-11
Salinity	0.5	0.5	ppt	1	02/14/20	AP	SM2520B-10
Mercury Digestion	Completed				02/17/20	LS/LS	SW7470A
Total Metals Digestion	Completed				02/14/20	AG	

Project ID: CHARLES RIVER Phoenix I.D.: CF32691

Client ID: CHARLES SW

RL/

Parameter Result PQL Units Dilution Date/Time By Reference

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

Ammonia:

This sample was received with a pH>2 The EPA requires preservation at time of sampling to a pH of <2. A sample bias can not be ruled out.

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

February 18, 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

February 18, 2020

QA/QC Data

SDG I.D.: GCF32691 % Sample Dup Dup LCS **LCSD** LCS MS MSD MS Rec RPD Blank **RPD RPD** Limits Limits RΙ Result Result **RPD** % % % % Parameter QA/QC Batch 518741 (mg/L), QC Sample No: CF32303 (CF32691) BRL 0.0002 <0.0002 <0.0002 NC Mercury - Water 117 120 75 - 125 30 Comment: Additional Mercury criteria: LCS acceptance range for waters is 80-120% and for soils is 75-125% QA/QC Batch 518646 (mg/L), QC Sample No: CF32121 (CF32691) ICP Metals - Aqueous Antimony BRL 0.005 < 0.005 < 0.005 NC 99.0 99.1 0.1 98.7 98.9 0.2 80 - 120 20 **BRL** 0.004 < 0.004 < 0.004 NC 95.1 95.5 95.6 Arsenic 0.4 96.5 0.9 80 - 120 20 Cadmium BRL 0.001 < 0.001 < 0.001 NC 95.9 97.2 1.3 95.4 95.9 0.5 80 - 120 20 0.001 Chromium BRL 0.001 0.001 NC 95.9 97.0 1.1 96.0 96.6 0.6 80 - 120 20 Copper BRL 0.005 < 0.005 < 0.005 NC 99.6 100 0.4 101 101 0.0 80 - 120 20 Iron BRL 0.010 0.072 0.073 1.40 97.5 99.1 1.6 97.7 98.4 80 - 120 20 91.8 BRL 0.002 < 0.002 < 0.002 NC 90.9 1.0 90.4 91.4 Lead 1.1 80 - 120 20 BRL 0.001 0.002 0.001 NC 95.9 97.0 1.1 95.6 96.0 80 - 120 Nickel 20 < 0.010 NC 94.7 80 - 120 Selenium **BRL** 0.010 < 0.010 93.7 1.1 94.4 94.8 0.4 20 Silver **BRL** 0.001 < 0.001 < 0.001 NC 98.7 99.8 1.1 99.8 101 1.2 80 - 120 20 80 - 120 Zinc BRL 0.004 0.062 0.062 0 96.0 97.0 1.0 96.6 97.0 0.4 20



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

QA/QC Report

February 18, 2020

QA/QC Data

SDG I.D.: GCF32691

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 518705 (PH), QC	Sample	No: Cl	32686 (0	CF32691)									
pH at 25C - Soil Comment:			7.92	7.90	0.30	100						85 - 115	20	
Additional: LCS acceptance range	e is 85-11	5% MS	acceptance	e range 7	75-125%									
QA/QC Batch 518677 (mg/L), C	2C Samp	ole No:	CF32635	(CF326	91)									
Chromium, Hexavalent Comment:	BRL	0.01	<0.01	<0.01	NC	101			114			90 - 110	30	
Additional Hexavalent Chromium	Additional Hexavalent Chromium criteria: LCS acceptance range for waters is 90-110% and MS acceptance range is 85-115%.													
QA/QC Batch 518644 (mg/L), C	2C Samp	ole No:	CF31644	(CF326	91)									
Ammonia as Nitrogen Comment:	BRL	0.05	25.6	25.0	2.40	100			107			90 - 110	20	
TKN is reported as Organic Nitrogen in the Blank, LCS, DUP and MS.														

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

February 18, 2020

Tuesday, February 18, 2020 Criteria: MA: CAM, GW1

Sample Criteria Exceedances Report GCF32691 - STANTECMA

State: MA

RLAnalysis SampNo Acode Phoenix Analyte Criteria Result RL Criteria 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.

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

MassDEP Analytical Protocol Certification Form											
Labo	Laboratory Name: Phoenix Environmental Laboratories, Inc. Project #:										
Proje	Project Location: CHARLES RIVER RTN:										
This F	orm provid	les certification	ons for	the following	data set	: [list Laborato	ry San	nple ID Number((s)]		
CF326	91										
Matric	es: 🗸 Gro	undwater/Sur	face Wa	ter Soil/S	Sedimen	t Drinking	y Wate	r 🗌 Air	Ot	her:	
CAM	Protocol (check all th	at app	y below)							
8260 V CAM II		7470/7471 H CAM III B	g •	MassDEP VPI CAM IV A	H	8081 Pesticides CAM V B		7196 Hex Cr CAM VI B		MassD CAM IX	EP APH
8270 S CAM II		7010 Metals CAM III C		MassDEP EPI CAM IV B	H	8151 Herbicides CAM V C		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	е		
	Affirmat	ive respons	es to q	uestions A t	hrough	F are require	d for	"Presumptive	Certai	nty" s	tatus
Α	Chain-of-Glaboratory	Custody, pro v, and prepar	perly p ed/ana	reserved (inc lyzed with me	luding to	nt with those d emperature*) ir olding times? (the fi see r	ield or narrative)	✓ \	Yes	□No
В		analytical m CAM protoco			ociated	QC requireme	nts sp	ecified in the	\	Yes	□ No
С		CAM protoco				I response acti htified performa			✓ \	Yes	□ No
D	CAM VII A		ssuran	e and Quality		orting requirem of Guidelines fo			>	Yes	□No
Е		t modification				method condu method(s) for				Yes	□No
			ethods	only: Was the	e compl	ete analyte list	repor	ted for each		Yes	□ No
F	Were all applicable CAM protocol QC and performance standard non-conformances identified and evaluated in a laboratory narrative (including all "No" responses to Questions A through E)? ✓ Yes □ No										
	Res	oonses to q	uestio	ns G, H and I	below	is required fo	r "Pre	sumptive Cer	tainty"	statu	s
G		reporting lim		r below all CA	AM repo	orting limits spe	cified	in the	✓ \	Yes	□ No
Data User Note: Data that achieve "Presumptive Certainty" status may not necessarily meet the data usability and representativeness requirements described in 310 CMR 40. 1056(2)(k) and WSC-07-350											
Н	Were all (QC performa	nce sta	ındards speci	fied in t	he CAM protoc	col(s) a	achieved?	~ \	Yes	□ No
l	Were resu protocol(s	•	for the	complete and	alyte lis	t specified in th	e sele	ected CAM	`	Yes	☑ No
1 41-			•	•				boratory narrativ			
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.											
						D	ate: ٦	Гuesday, Febi	ruary 1	18, 20	20
	orized	RN	١٨٠٠.	: NWK	~	Printed Na	me: F	Rashmi Makol	- 		
Sign	Authorized Signature: Printed Name: Rashmi Makol Position: Project Manager										



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



MCP Certification Report

February 18, 2020 SDG I.D.: GCF32691

SDG Comments

Metals Analysis:

The client requested a site specific list of elements which is shorter than the 6010 MCP list.

Hexavalent Chromium (Aqueous)

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

Instrument:

BECKMAN DU720 02/14/20-1 Dustin Harrison, Chemist 02/14/20

CF32691

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

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

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

QC (Batch Specific):

Batch 518677 (CF32635)

CF32691

All LCS recoveries were within 90 - 110 with the following exceptions: None.

Additional Hexavalent Chromium criteria: LCS acceptance range for waters is 90-110% and MS acceptance range is 85-115%.

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 02/17/20 09:49 Rick Schweitzer, Chemist 02/17/20

CF32691

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 518741 (CF32303)

CF32691

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%



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



Certification Report

February 18, 2020 SDG I.D.: GCF32691

ICP Metals Narration

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

Instrument:

BLUE 02/15/20 08:02

Cindy Pearce, Chemist 02/15/20

CF32691

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 518646 (CF32121)

CF32691

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.

NITROGEN

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

Instrument:

LACHAT 02/15/20-1

Kandi Della Bella, Chemist 02/15/20

CF32691

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

All method verification standards and blanks met criteria.

QC (Batch Specific):

Batch 518644 (CF31644)

CF32691

All LCS recoveries were within 85 - 115 with the following exceptions: None.

TKN is reported as Organic Nitrogen in the Blank, LCS, DUP and MS.

Additional criteria: LCS acceptance range for waters is 85-115% and for soils is 75-125%. MS acceptance range is 75-125%.

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.

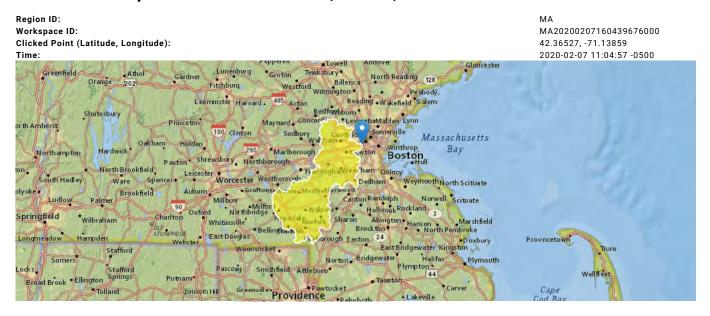
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CHAIN OF CLISTODY RECORD		East Middle Turnpike, P.O. Box 370, Manchester, CT 06040 Email: info@phoenixlabs.com Fax (860) 645-0823 Client Services (860) 645-8726	Creyes	300		Richard CI MA
CHAIN		587 East Middle Turnpike, P.O. Box Email: info@phoenixlabs.com Client Services (8	Project:	Report to Invoice to QUOTE#	A A A A A A A A A A A A A A A A A A A	Date: Time: 2.17.23 5.15 Turnaround Time: 1.04 1.08 2.08 2.08 3.08 3.08 0.04 0.05 0.04 0.04 0.04 0.04 0.05
		es, Inc.	W + 1/5	Colony Prov	Surface Water www=wasteseoil SD=Soil w=wipe Sample Date Matrix Sampled Sw 2442b	by: Home: X 1 days Silve, Iran, Silve, Zinc,
		Environmental Laboratories, Inc.	STANKC CONSULTION	400 count	Sampler's Signature Signature Matrix Code: DW=Drinking Water GW=Ground Water SW=Surface Water WW=Waste Water RW=Raw Water SE=Sediment SL=Sludge S=Soil SD=Soild W=Wipe OIL=Oil B=Bulk L=Liquid X = (Other) Customer Sample Sample Date Time Identification Matrix Sampled Sample SAMPLE # (Other) SAMPLE # (DW-CK)	Relinquished by: Relinquished
	DI IC	Environme	Customer:	Address:	Sampler's Signature Matrix Code: DW=Drinking Water RW=Raw Water St B=Bulk L=Liquid X PHOENIX USE ONLY SAMPLE #	Relinguished by: Comments, Special Anthwarly, Assetting Mercury, I



APPENDIX C
STREAMSTATS 7Q10 FLOW REPORT, DILUTION FACTOR CALCULATIONS AND COPY OF MALIMITSBOOKREV1.XLSX

StreamStats Page 1 of 2

StreamStats Report - 30 Penniman Road, Allston, MA



Basin Characteristics			
Parameter Code	Parameter Description	Value	Unit
DRNAREA	Area that drains to a point on a stream	279	square miles
BSLDEM250	Mean basin slope computed from 1:250K DEM	2.342	percent
DRFTPERSTR	Area of stratified drift per unit of stream length	0.23	square mile per mile
MAREGION	Region of Massachusetts 0 for Eastern 1 for Western	0	dimensionless

Low-Flow Statistics Parameters[Statewide Low Flow WRIR00 4135]						
Parameter Code	Parameter Name	Value	Units	Min Limit	Max Limit	
DRNAREA	Drainage Area	279	square miles	1.61	149	
BSLDEM250	Mean Basin Slope from 250K DEM	2.342	percent	0.32	24.6	
DRFTPERSTR	Stratified Drift per Stream Length	0.23	square mile per mile	0	1.29	
MAREGION	Massachusetts Region	0	dimensionless	0	1	

StreamStats Page 2 of 2

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

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

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

Statistic	Value	Unit
7 Day 2 Year Low Flow	48.9	ft^3/s
7 Day 10 Year Low Flow	24.3	ft^3/s

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/)

USGS Data Disclaimer: Unless otherwise stated, all data, metadata and related materials are considered to satisfy the quality standards relative to the purpose for which the data were collected. Although these data and associated metadata have been reviewed for accuracy and completeness and approved for release by the U.S. Geological Survey (USGS), no warranty expressed or implied is made regarding the display or utility of the data for other purposes, nor on all computer systems, nor shall the act of distribution constitute any such warranty.

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Application Version: 4.3.11

STANTEC CONSULTING SERVICES, INC.

30 Penn, LLC

CALCULATIONS

PROJECT NO. 195601892

SHEET

1 of

2020 February

1

DATE

COMPUTED BY

RBL

PROJECT SUBJECT

CLIENT

30 Penniman Road Redevelopment **Dilution Factor Calculations**

PURPOSE:

Calculate Dilution Factor (DF) for project based on 7 Day 10 Year (7Q10) Low Flow values.

APPROACH:

Calculate DF based on EPA formula $(Q_S + Q_D)/Q_D$, where Q_S is 7Q10 in million gallons per day (MGD) and Q_D is discharge flow

in MGD.

ASSUMPTIONS: 1. 7Q10 is 24.3 cfs (from StreamStats 4.0)

2. A conversion of 7.48 is used to convert cubic feet to gallons

3. A discharge flowrate of 500 gpm is assumed

CALCULATIONS:

7Q10 Low Flow Value (Q_s)

24.3 ft³ Χ $Q_S =$ sec

7.48 gallons

8<u>6,400 sec</u> Χ day

Χ

<u> 1 MG</u> 1,000,000

gallons

 $Q_s =$

15.70 MGD

Discharge Flowrate (Q_D)

500 gallons $Q_D =$ min

Χ

1,440 min day

<u>1 MG</u> 1,000,000 gallons

22.81

 $Q_D = 0.72 \text{ MGD}$

Dilution Factor (DF)

DF =

 $Q_s + Q_D$ 15.70 MGD + 0.72 MGD 0.72 MGD

CONCLUSION

The dilution factor for this project is calculated to be 22.81 based on the provided 7Q10 low flow value and discharge flowrate.

RE: Allston RGP 7Q10 and Dilution Factor

Vakalopoulos, Catherine (DEP) <catherine.vakalopoulos@state.ma.us>

Wed 2/12/2020 3:31 PM

To: Learned, Richard <Richard.Learned@stantec.com>
Cc: Salvetti, Joseph <Joseph.Salvetti@stantec.com>

Hi Rich.

Yes, nice speaking with you today. I can confirm that your dilution factor calculation (22.81) for this proposed discharge from 30 Penniman Rd. in Allston to a storm drain that discharges to the Charles River near the Artesani Playground is correct.

To assist you with filling out the NOI, this segment of the Charles River is identified as MA72-36, is classified as Class B, is not an Outstanding Resource Water, and there are two approved TMDLs for pathogens and phosphorus. To see the causes of impairments, go to:

https://www.mass.gov/doc/final-massachusetts-year-2016-integrated-list-of-waters/download and search for "MA72-36".

As we discussed, since this is a *current* MCP site, you do not have to submit a transmittal form and \$500 fee (https://www.mass.gov/how-to/wm-15-npdes-general-permit-notice-of-intent). You can come on the NOI but please don't send me a paper copy.

Feel free to contact me if you have any additional questions.

Cathy

Cathy Vakalopoulos, Massachusetts Department of Environmental Protection 1 Winter St., Boston, MA 02108, 617-348-4026

A Please consider the environment before printing this e-mail

From: Learned, Richard [mailto:Richard.Learned@stantec.com]

Sent: Wednesday, February 12, 2020 2:50 PM

To: Vakalopoulos, Catherine (DEP)

Cc: Salvetti, Joseph

Subject: Allston RGP 7Q10 and Dilution Factor

Cathy,

It was a pleasure to speak with you today.

Attached please find the StreamStats 7Q10 flow report and dilution calcs for a dewatering project we are planning in Allston, MA. The site is located at 30 Penniman Road. Excavation below the water table is necessary to construct the building foundation. We are using extraction wells located around the excavation to dewater. Groundwater from the wells will be pumped into a Frac tank, through filtration and through granular activated carbon (GAC) before discharge into a BWSC stormwater catch basin in Penniman Road (#172). The stormwater system discharges into the Charles River at outfall SDO037 (see plan).

We are using a maximum flow rate of 500 gpm with an expected average rate of 250 gpm.

Please verify our 7Q10 flow and dilution calcs consistent with Appendix V of the RGP NOI.

Please let me know if you need any other information.

Richard Learned LSP

Senior Environmental Project Manager

Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stantec 146 Main Street Unit 3 Hyannis MA 02601-3128 US

Stantec

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Enter number values in green boxes based on the instructions to the right

Enter values in the units specified

15.7 Q_R = Enter upstream flow in MGD 0.36 Q_P = Enter discharge flow in MGD 0 Downstream 7Q10

Enter a dilution factor for saltwater receiving water (this box does not apply to freshwater receiving waters)



Enter values in the units specified

↓
 567 C_d = Enter influent hardness in mg/L CaCO₃
 74.4 C_s = Enter receiving water hardness in mg/L CaCO₃

Enter receiving water concentrations in the units specified

\downarrow	_	Impaired for metals?
7.31	pH in Standard Units	\downarrow
12	Temperature in °C	
0.14	Ammonia in mg/L	
74.4	Hardness in mg/L CaCO	3
0.5	Salinity in ppt	
<5	Antimony in μg/L	no
<4	Arsenic in μg/L	no
<1	Cadmium in µg/L	yes
<10	Chromium III in µg/L	yes
<10	Chromium VI in μg/L	yes
<5	Copper in µg/L	yes
494	Iron in μg/L	yes
<2	Lead in μg/L	yes
< 0.2	Mercury in μg/L	yes
<1	Nickel in μg/L	yes
<10	Selenium in μg/L	yes
<1	Silver in μg/L	yes
<9	Zinc in μg/L	yes

Enter influent concentrations in the units specified

\downarrow	
0	TRC in μg/L
0.7	Ammonia in mg/L
0	Antimony in μg/L
11	Arsenic in μg/L
0	Cadmium in µg/L
9	Chromium III in µg/L
0	Chromium VI in µg/L
5	Copper in µg/L
10000	Iron in μg/L
9	Lead in μg/L
0	Mercury in μg/L
11	Nickel in μg/L
0	Selenium in μg/L
0	Silver in μg/L
34	Zinc in μg/L
0	Cyanide in µg/L
0	Phenol in μg/L
0	Carbon Tetrachloride in µg/L
100	Tetrachloroethylene in μg/L
0	Total Phthalates in μg/L
0	Diethylhexylphthalate in μg/L
0	Benzo(a)anthracene in μg/L
0	Benzo(a)pyrene in μg/L
0	Benzo(b)fluoranthene in μg/L
0	Benzo(k)fluoranthene in μg/L
0	Chrysene in μg/L
0	Dibenzo(a,h)anthracene in μg/L
0	Indeno(1,2,3-cd)pyrene in μg/L
0	Methyl-tert butyl ether in μg/L

Notes: Revised 1-24-20

Freshwater: leave 0 unless 7Q10 or alternate Q_R <u>AND</u> a dilution factor >1 approved by the State; Saltwater (estuarine and marine): leave 0 unless QR approved by the State Enter the design flow or 1 MGD, whichever is less (100 gpm design flow = 0.144 MGD and is entered by default) Leave 0 unless Q_R approved by the State

Freshwater: leave 0

Saltwater (estuarine and marine): leave 0 unless DF approved by the State

Applies to freshwater receiving waters only

pH, temperature, and ammonia required for all discharges
Hardness required for freshwater
Salinity required for saltwater (estuarine and marine)
Metals required for all discharges if detected in the influent and if dilution factor approved by State
Enter 0 if non-detect or testing not required

If receiving water is not listed as impaired for metals in State 303(d) List, change to "no" using dropdown

if >1 sample, enter maximum influent measurement

if>10 samples, may enter 95th percentile of influent measurements using EPA's Technical Support Document for Water Quality-based Toxics Control

Enter 0 if non-detect or testing not required



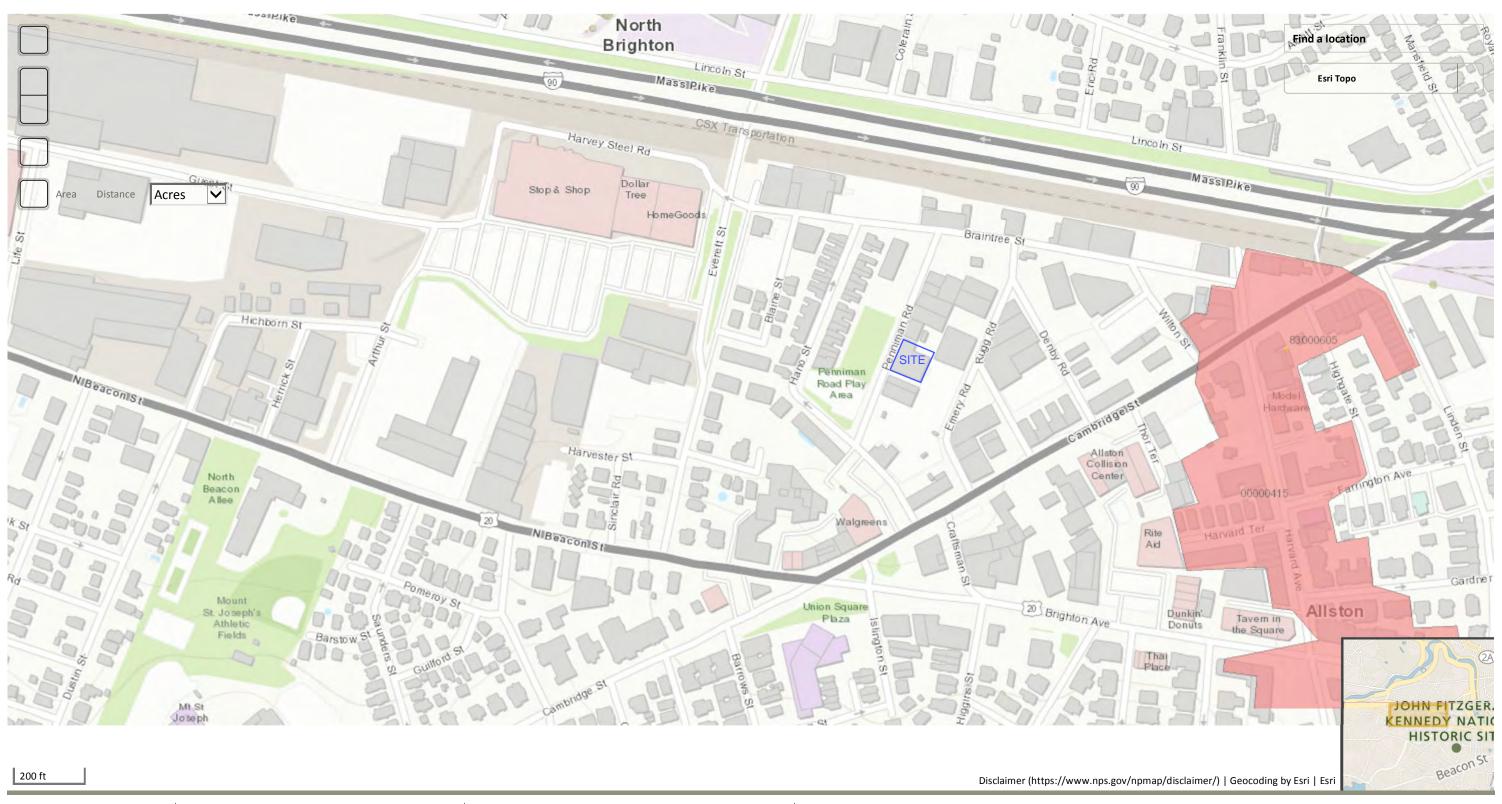
APPENDIX D
NATIONAL REGISTER OF HISTORIC PLACES DOCUMENTATION

National Register of Historic Places

National Register of Historic Places

National Park Service
U.S. Department of the Interior

Public, non-restricted data depicting National Register spatial data processed by the Cultural Resources GIS facility. Data last updated in April, 2014.



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Massachusetts Cultural Resource Information System MACRIS

MACRIS Search Results

Search Criteria: Town(s): Boston; Place: Allston; Street No: 30; Street Name: PENNIMAN; Resource Type(s): Building;

Inv. No. Property Name Street Town Year

Friday, February 7, 2020 Page 1 of 1



APPENDIX E
ENDANGERED SPECIES ACT DOCUMENTATION

IPaC: Explore Location Page 1 of 12

IPaC Information for Planning and Consultation 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

Suffolk County, Massachusetts



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

IPaC: Explore Location Page 2 of 12

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.
- Click REQUEST SPECIES LIST.

Listed species

¹ 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>.

- 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.

THERE ARE NO ENDANGERED SPECIES EXPECTED TO OCCUR AT THIS LOCATION.

IPaC: Explore Location Page 3 of 12

Migratory birds

Certain birds are protected under the Migratory Bird Treaty Act

¹ and the Bald and Golden Eagle Protection Act².

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 <u>Bald and Golden Eagle Protection Act</u> 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>.

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"

IPaC: Explore Location Page 4 of 12

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

Breeds Oct 15 to Aug 31

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

Breeds May 15 to Oct 10

Bobolink Dolichonyx oryzivorus

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

Breeds May 20 to Jul 31

Canada Warbler Cardellina canadensis

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

Breeds May 20 to Aug 10

Cerulean Warbler Dendroica cerulea

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

https://ecos.fws.gov/ecp/species/2974

Breeds Apr 29 to Jul 20

Dunlin Calidris alpina arcticola

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

Breeds elsewhere

Evening Grosbeak Coccothraustes vespertinus

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

Breeds elsewhere

Kentucky Warbler Oporornis formosus

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

Breeds Apr 20 to Aug 20

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

IPaC: Explore Location Page 5 of 12

Nelson's Sparrow Ammodramus nelsoni

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

Breeds May 15 to Sep 5

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

Prothonotary Warbler Protonotaria citrea

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

Breeds Apr 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

Red-throated Loon Gavia stellata

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

Breeds elsewhere

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

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 (■)

IPaC: Explore Location

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.
- 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.

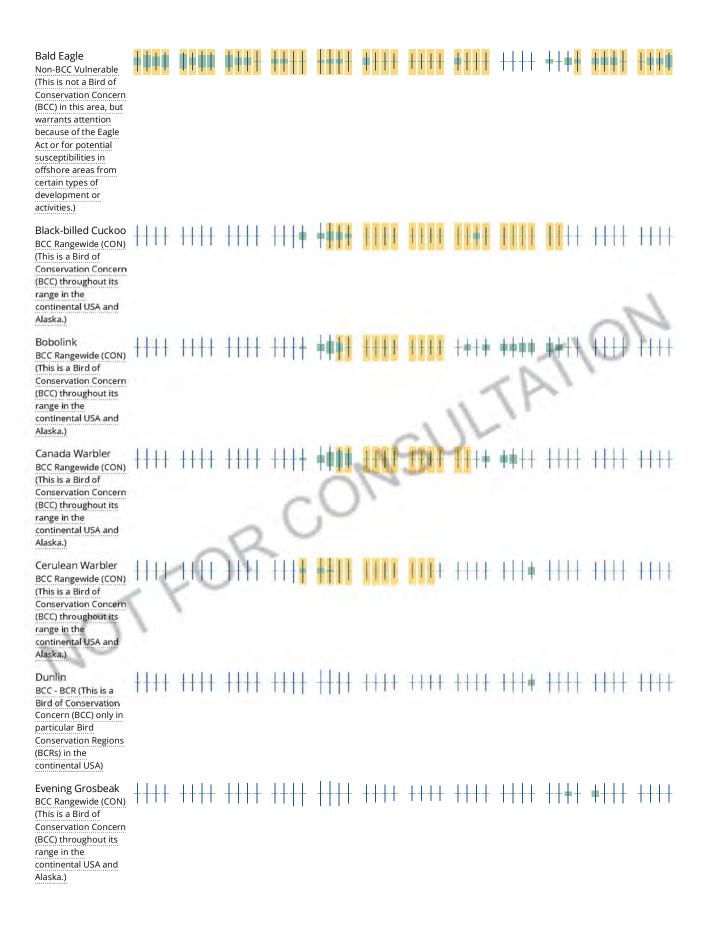
No Data (-)

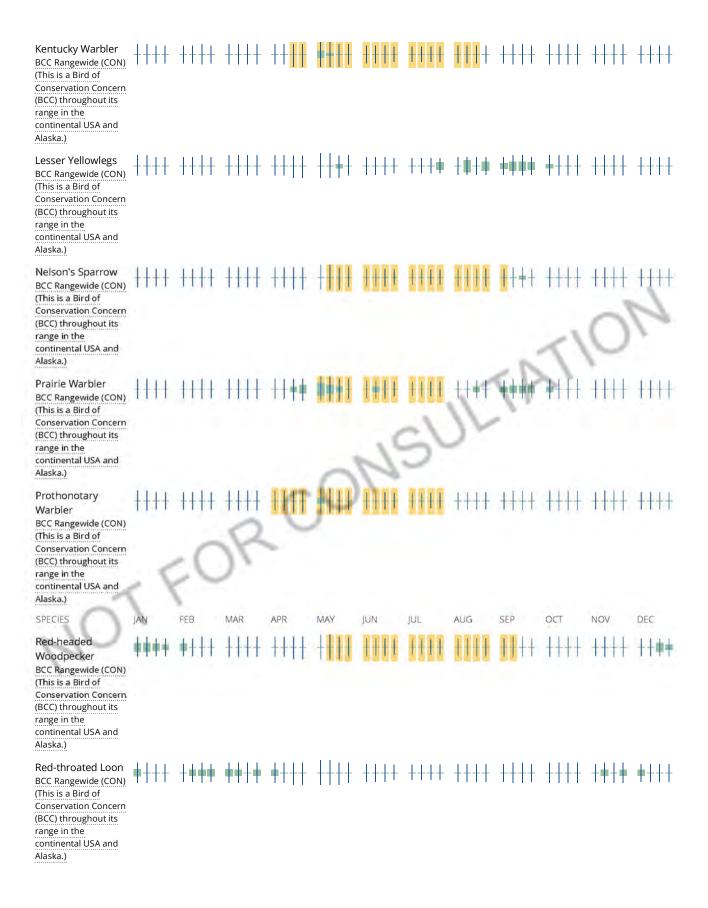
A week is marked as having no data if there were no survey events for that week.

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</u>, <u>banding</u>, <u>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 <u>AKN Phenology Tool</u>.

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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 science datasets</u>.

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:

- "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
- "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

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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.

Facilities

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</u> <u>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:

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RIVERINE

R2UBH

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

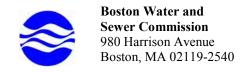
Certain wetland habitats are excluded from the National mapping program because of the limitations of aerial imagery as the primary data source used to detect wetlands. These habitats include seagrasses or submerged aquatic vegetation that are found in the intertidal and subtidal zones of estuaries and 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.



APPENDIX F
COPIES OF BWSC PERMIT APPLICATIONS AND DCR CORRESPONDENCE



DEWATERING DISCHARGE PERMIT APPLICATION

OWNER / AUTHORIZED APPLICANT PROVIDE INFORMATION HERE:

Company Name:	Address:
Phone Number:	Fax number:
Contact person name:	Title:
Cell number:	Email address:
	ew Application Permit Extension Other (Specify):
Owner's Information (if different	from above):
Owner of property being dewatered	l:
	Phone number:
Location of Discharge & Propose	d Treatment System(s):
Street number and name:	Neighborhood
Discharge is to a: ☐ Sanitary Sewe	er Combined Sewer Storm Drain Other (specify):
Describe Proposed Pre-Treatment S	System(s):
	Receiving Waters
	nticipated Dates of Discharge): FromTo □ Tank Removal/Installation □ Foundation Excavation
☐ Groundwater Remediation☐ Utility/Manhole Pumping	☐ Tank Removal/Installation ☐ Foundation Excavation ☐ Test Pipe ☐ Trench Excavation
□ Accumulated Surface Water	☐ Hydrogeologic Testing ☐ Other
Permanent Discharges	
□ Foundation Drainage	□ Crawl Space/Footing Drain
□ Accumulated Surface Water□ Non-contact/Uncontaminated Proces	□ Non-contact/Uncontaminated Cooling □ Other;
number, size, make and start reading. 1. If discharging to a sanitary or combine as other relevant information.	of the discharge and the location of the point of discharge (i.e. the sewer pipe or catch basin). Include meter type, meter Note. All discharges to the Commission's sewer system will be assessed current sewer charges. In discharge to the Commission's sewer system will be assessed current sewer charges. In discharge, attach a copy of MWRA's Sewer Use Discharge permit or application. In attach a copy of EPA's NPDES Permit or NOI application, or NPDES Permit exclusion letter for the discharge, as well senied or revoked if applicant fails to obtain the necessary permits from MWRA or EPA. Boston Water and Sewer Commission Engineering Customer Services 980 Harrison Avenue, Boston, MA 02119 Attn: Jodi Dobay, Engineering Customer Service E-mail: beginj@bwsc.org Phone: 617-989-7259 Fax: 617-989-7716

RE: Allston Outfall

Casey, Sean (DCR) <sean.casey@state.ma.us>

Wed 2/12/2020 9:58 AM

To: Learned, Richard < Richard.Learned@stantec.com>

Thanks Rich. It is confirmed that this is not a DCR asset and does not require DCR permitting for its use.

Thanks, Sean

From: Learned, Richard < Richard.Learned@stantec.com>

Sent: Wednesday, February 12, 2020 9:41 AM **To:** Casey, Sean (DCR) <Sean.Casey@mass.gov>

Subject: Re: Allston Outfall

Sean,

See attached.

Richard Learned LSP

Senior Environmental Project Manager

Direct: 508 591-4351 Cell: 508 326-9913

Richard.Learned@stantec.com

Stantec 146 Main Street Unit 3 Hyannis MA 02601-3128 US

Stantec

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From: Casey, Sean (DCR) < sean.casey@state.ma.us > Sent: Wednesday, February 12, 2020 9:39 AM

To: Learned, Richard < Richard.Learned@stantec.com

Subject: Allston Outfall

Rich,

Thank you for your voicemail and I have inquired with the DCR Stormwater team. Unfortunately the number you provided is not how DCR labels their outfalls. They have asked if you could send a

map/plan/google image, so that we can better determine if the outfall is DCR's and whether you will need a permit.

Thanks,

SEAN CASEY
Director- Construction/Access Permits & Accident Recovery
Department of Conservation and Recreation
251 Causeway St, Suite 700
Boston, MA 02114
617-626-1444



APPENDIX G BEST MANAGEMENT PRACTICES PLAN (BMPP)

Consistent with Section 2.5(2)(d) of the NPDES NOI, a Pollutant Minimization BMP has been prepared for the proposed discharge. The BMP includes:

- i. Identification and assessment of the type and quantity of pollutants, including their potential to impact receiving water quality;
- ii. Water quality control measures that ensure dilution is not used as a form of treatment, or as a means to achieve the limitations and requirements in the RGP general permit; and
- iii. Selection, design, installation and proper operation and maintenance of pollution control technologies necessary to meet the limitations and requirements in the RGP general permit. The treatment technologies may include, but are not limited to any combination of the following:
 - 1) Adsorption/Absorption
 - 2) Granulated Activated Carbon (GAC)/Liquid Phase Carbon Adsorption
 - 3) Precipitation/Coagulation/Flocculation
 - 4) Separation/Filtration

The BMP will be available on-Site during the period of dewatering.