Site Management Periodic Review Report and IC/EC Certification (2022) Geneseo -Park Street MGP Site (V00731) Geneseo, New York December 2022

Exhibit A

Laboratory Reports for Groundwater Sampling Events

(Presented as separate file)



Analytical Report For

Neu-Velle

For Lab Project ID

214958

Referencing

RGE Geneseo Fmr. MGP Site

Prepared

Tuesday, November 16, 2021

Any noncompliant QC parameters or other notes impacting data interpretation are flagged or documented on the final report or are noted below:

Portions of the enclosed report reflects analysis that has been subcontracted and are presented in their original form.

A complete ASP package will follow this report.

Certifies that this report has been approved by the Technical Director or Designee

179 Lake Avenue • Rochester, NY 14608 • (585) 647-2530 • Fax (585) 647-3311 • ELAP ID# 10958



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW4-102921

Lab Sample ID:214958-01Date Sampled:10/29/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Anal	yzed
Benzene	< 1.00	ug/L		M	11/3/2021	13:33
Ethylbenzene	< 2.00	ug/L		M	11/3/2021	13:33
m,p-Xylene	2.44	ug/L			11/3/2021	13:33
o-Xylene	< 2.00	ug/L			11/3/2021	13:33
Toluene	< 2.00	ug/L		M	11/3/2021	13:33
Surrogate	Percei	nt Recovery	<u>Limits</u>	<u>Outliers</u>	Date Analy	zed
1,2-Dichloroethane-d4		121	77.9 - 132		11/3/2021	13:33
4-Bromofluorobenzene		118	62.6 - 133		11/3/2021	13:33
Pentafluorobenzene		115	88.9 - 114	*	11/3/2021	13:33
Toluene-D8		107	75.6 - 117		11/3/2021	13:33

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05208.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW1-103021

Lab Sample ID:214958-02Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>vzed</u>
Benzene	< 1.00	ug/L			11/3/2021	14:31
Ethylbenzene	< 2.00	ug/L			11/3/2021	14:31
m,p-Xylene	< 2.00	ug/L			11/3/2021	14:31
o-Xylene	< 2.00	ug/L			11/3/2021	14:31
Toluene	< 2.00	ug/L			11/3/2021	14:31
<u>Surrogate</u>	Percei	<u>ıt Recovery</u>	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		117	77.9 - 132		11/3/2021	14:31
4-Bromofluorobenzene		108	62.6 - 133		11/3/2021	14:31
Pentafluorobenzene		117	88.9 - 114	*	11/3/2021	14:31
Toluene-D8		108	75.6 - 117		11/3/2021	14:31

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05211.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW2-103021

Lab Sample ID:214958-03Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>vzed</u>
Benzene	< 1.00	ug/L			11/3/2021	14:50
Ethylbenzene	< 2.00	ug/L			11/3/2021	14:50
m,p-Xylene	1.10	ug/L		J	11/3/2021	14:50
o-Xylene	< 2.00	ug/L			11/3/2021	14:50
Toluene	< 2.00	ug/L			11/3/2021	14:50
Surrogate	Percei	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		118	77.9 - 132		11/3/2021	14:50
4-Bromofluorobenzene		107	62.6 - 133		11/3/2021	14:50
Pentafluorobenzene		113	88.9 - 114		11/3/2021	14:50
Toluene-D8		107	75.6 - 117		11/3/2021	14:50

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05212.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-DUP-103021

Lab Sample ID:214958-04Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>vzed</u>
Benzene	< 1.00	ug/L			11/3/2021	15:09
Ethylbenzene	< 2.00	ug/L			11/3/2021	15:09
m,p-Xylene	< 2.00	ug/L			11/3/2021	15:09
o-Xylene	< 2.00	ug/L			11/3/2021	15:09
Toluene	< 2.00	ug/L			11/3/2021	15:09
Surrogate	Percei	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		123	77.9 - 132		11/3/2021	15:09
4-Bromofluorobenzene		106	62.6 - 133		11/3/2021	15:09
Pentafluorobenzene		114	88.9 - 114		11/3/2021	15:09
Toluene-D8		111	75.6 - 117		11/3/2021	15:09

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05213.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW3-103121

Lab Sample ID:214958-05Date Sampled:10/31/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>yzed</u>
Benzene	< 1.00	ug/L			11/3/2021	15:29
Ethylbenzene	< 2.00	ug/L			11/3/2021	15:29
m,p-Xylene	< 2.00	ug/L			11/3/2021	15:29
o-Xylene	< 2.00	ug/L			11/3/2021	15:29
Toluene	< 2.00	ug/L			11/3/2021	15:29
<u>Surrogate</u>	<u>Percer</u>	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		120	77.9 - 132		11/3/2021	15:29
4-Bromofluorobenzene		118	62.6 - 133		11/3/2021	15:29
Pentafluorobenzene		110	88.9 - 114		11/3/2021	15:29
Toluene-D8		107	75.6 - 117		11/3/2021	15:29

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05214.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW6-103121

Lab Sample ID:214958-06Date Sampled:10/31/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>yzed</u>
Benzene	178	ug/L			11/8/2021	15:52
Ethylbenzene	32.6	ug/L			11/8/2021	15:52
m,p-Xylene	43.5	ug/L			11/8/2021	15:52
o-Xylene	43.3	ug/L			11/8/2021	15:52
Toluene	33.8	ug/L			11/8/2021	15:52
<u>Surrogate</u>	Perc	cent Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		99.0	77.9 - 132		11/8/2021	15:52
4-Bromofluorobenzene		86.5	62.6 - 133		11/8/2021	15:52
Pentafluorobenzene		99.3	88.9 - 114		11/8/2021	15:52
Toluene-D8		88.7	75.6 - 117		11/8/2021	15:52

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05290.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-EB-110121

Lab Sample ID:214958-07Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>vzed</u>
Benzene	0.806	ug/L		J	11/3/2021	16:07
Ethylbenzene	< 2.00	ug/L			11/3/2021	16:07
m,p-Xylene	< 2.00	ug/L			11/3/2021	16:07
o-Xylene	< 2.00	ug/L			11/3/2021	16:07
Toluene	< 2.00	ug/L			11/3/2021	16:07
Surrogate	Percei	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		117	77.9 - 132		11/3/2021	16:07
4-Bromofluorobenzene		117	62.6 - 133		11/3/2021	16:07
Pentafluorobenzene		116	88.9 - 114	*	11/3/2021	16:07
Toluene-D8		108	75.6 - 117		11/3/2021	16:07

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05216.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW8-110121

Lab Sample ID:214958-08Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	Res	<u>Sult</u> <u>Units</u>		Qualifier	Date Analy	vzed
Benzene	3.37	ug/L			11/3/2021	16:27
Ethylbenzene	2.32	ug/L			11/3/2021	16:27
m,p-Xylene	3.52	ug/L			11/3/2021	16:27
o-Xylene	3.06	ug/L			11/3/2021	16:27
Toluene	1.44	ug/L		J	11/3/2021	16:27
<u>Surrogate</u>		Percent Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		118	77.9 - 132		11/3/2021	16:27
4-Bromofluorobenzene		112	62.6 - 133		11/3/2021	16:27
Pentafluorobenzene		115	88.9 - 114	*	11/3/2021	16:27
Toluene-D8		97.0	75.6 - 117		11/3/2021	16:27

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05217.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW7-110121

Lab Sample ID:214958-09Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Analy	<u>vzed</u>
Benzene	< 1.00	ug/L			11/3/2021	16:46
Ethylbenzene	< 2.00	ug/L			11/3/2021	16:46
m,p-Xylene	< 2.00	ug/L			11/3/2021	16:46
o-Xylene	< 2.00	ug/L			11/3/2021	16:46
Toluene	< 2.00	ug/L			11/3/2021	16:46
<u>Surrogate</u>	Percei	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		115	77.9 - 132		11/3/2021	16:46
4-Bromofluorobenzene		101	62.6 - 133		11/3/2021	16:46
Pentafluorobenzene		114	88.9 - 114		11/3/2021	16:46
Toluene-D8		97.7	75.6 - 117		11/3/2021	16:46

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05218.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: Trip Blank T1075

Lab Sample ID:214958-10Date Sampled:10/25/2021Matrix:WaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Anal	<u>yzed</u>
Benzene	< 1.00	ug/L			11/3/2021	17:05
Ethylbenzene	< 2.00	ug/L			11/3/2021	17:05
m,p-Xylene	< 2.00	ug/L			11/3/2021	17:05
o-Xylene	< 2.00	ug/L			11/3/2021	17:05
Toluene	< 2.00	ug/L			11/3/2021	17:05
<u>Surrogate</u>	Percei	nt Recovery	<u>Limits</u>	Outliers	Date Analy	zed
1,2-Dichloroethane-d4		127	77.9 - 132		11/3/2021	17:05
4-Bromofluorobenzene		111	62.6 - 133		11/3/2021	17:05
Pentafluorobenzene		117	88.9 - 114	*	11/3/2021	17:05
Toluene-D8		121	75.6 - 117	*	11/3/2021	17:05

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05219.D



Analytical Report Appendix

The reported results relate only to the samples as they have been received by the laboratory.

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All soil/sludge samples have been reported on a dry weight basis, unless qualified "reported as received". Other solids are reported as received.

Low level Volatiles blank reports for soil/solid matrix are based on a nominal 5 gram weight. Sample results and reporting limits are based on actual weight, which may be more or less than 5 grams.

The Chain of Custody provides additional information, including compliance with sample condition requirements upon receipt. Sample condition requirements are defined under the 2003 NELAC Standard, sections 5.5.8.3.1 and 5.5.8.3.2.

NYSDOH ELAP does not certify for all parameters. Paradigm Environmental Services or the indicated subcontracted laboratory does hold certification for all analytes where certification is offered by ELAP unless otherwise specified. Aliquots separated for certain tests, such as TCLP, are indicated on the Chain of Custody and final reports with an "A" suffix.

Data qualifiers are used, when necessary, to provide additional information about the data. This information may be communicated as a flag or as text at the bottom of the report. Please refer to the following list of analyte-specific, frequently used data flags and their meaning:

- "<" = Analyzed for but not detected at or above the quantitation limit.
- "E" = Result has been estimated, calibration limit exceeded.
- "Z" = See case narrative.
- "D" = Sample, Laboratory Control Sample, or Matrix Spike Duplicate results above Relative Percent Difference limit.
- "M" = Matrix spike recoveries outside QC limits. Matrix bias indicated.
- "B" = Method blank contained trace levels of analyte. Refer to included method blank report.
- "I" = Result estimated between the quantitation limit and half the quantitation limit.
- "L" = Laboratory Control Sample recovery outside accepted QC limits.
- "P" = Concentration differs by more than 40% between the primary and secondary analytical columns.
- "NC" = Not calculable. Applicable to RPD if sample or duplicate result is non-detect or estimated (see primary report for data flags). Applicable to MS if sample is greater or equal to ten times the spike added. Applicable to sample surrogates or MS if sample dilution is 10x or higher.
- "*" = Indicates any recoveries outside associated acceptance windows. Surrogate outliers in samples are presumed matrix effects. LCS demonstrates method compliance unless otherwise noted.
- "(1)" = Indicates data from primary column used for QC calculation.
- "A" = denotes a parameter for which ELAP does not offer approval as part of their laboratory certification program.
- "F" = denotes a parameter for which Paradigm does not carry certification, the results for which should therefore only be used where ELAP certification is not required, such as personal exposure assessment.

This report is part of a multipage document and should only be evaluated in its entirety. The Chain of Custody provides additional sample information, including compliance with the sample condition requirements upon receipt.

GENERAL TERMS AND CONDITIONS LABORATORY SERVICES

These Terms and Conditions embody the whole agreement of the parties in the absence of a signed and executed contract between the Laboratory (LAB) and Client. They shall supersede all previous communications, representations, or agreements, either verbal or written, between the parties. The LAB specifically rejects all additional, inconsistent, or conflicting terms, whether printed or otherwise set forth in any purchase order or other communication from the Client to the LAB. The invalidity or unenforceability in whole or in part of any provision, tern or condition hereof shall not affect in any way the validity or enforceability of the remainder of the Terms and Conditions. No waiver by LAB of any provision, term, or condition hereof or of any breach by or obligation of the Client hereunder shall constitute a waiver of such provision, term, or condition on any other occasion or a waiver of any other breach by or obligation of the Client. This agreement shall be administered and interpreted under the laws of the state which services are procured.

Warranty.

Recognizing that the nature of many samples is unknown and that some may contain potentially hazardous components, LAB warrants only that it will perform testing services, obtain findings, and prepare reports in accordance with generally accepted analytical laboratory principles and practices at the time of performance of services. LAB makes no other warranty, express or implied.

Scope and Compensation. LAB agrees to perform the services described in the chain of custody to which these terms and conditions are attached. Unless the parties agree in writing to the contrary, the duties of LAB shall not be construed to exceed the services specifically described. LAB wi use LAB default method for all tests unless specified otherwise on the Work Order.

Payment terms are net 30 days from the date of invoice. All overdue payments are subject to an interest charge of one and one-half percent (1-1/2%) per month or a portion thereof. Client shall also be responsible for costs of collection, including payment of reasonable attorney fees if such expense is incurred. The prices, unless stated, do not include any sale, use or other taxes. Such taxes will be added to invoice prices when required.

Prices.

Compensation for services performed will be based on the current Lab Analytical Fee Schedule or on quotations agreed to in writing by the parties. Turnaround time based charges are determined from the time of resolution of all work order questions. Testimony, court appearances or data compilation for legal action will be charged separately. Evaluation and reporting of initial screening runs may incur additional fees.

Limitations of Liability.

In the event of any error, omission, or other professional negligence, the sole and exclusive responsibility of LAB shall be to reperform the deficient work at its own expense and LAB shall have no other liability whatsoever. All claims shall be deemed waived unless made in writing and received by LAB within ninety (90) days following completion of services.

LAB shall have no liability, obligation, or responsibility of any kind for losses, costs, expenses, or other damages (including but not limited to any special, direct, incidental or consequential damages) with respect to LAB's services or results.

All results provided by LAB are strictly for the use of its clients and LAB is in no way responsible for the use of such results by clients or third parties. All reports should be considered in their entirety, and LAB is not responsible for the separation, detachment, or other use of any portion of these reports. Client may not assign the lab report without the written consent of the LAB. Client covenants and agrees, at its/his/her sole expense, to indemnify, protect, defend, and save harmless the LAB from and against

any and all damages, losses, liabilities, obligations, penalties, claims, litigation, demands, defenses, judgments, suits, actions, proceedings, costs, disbursements and/or expenses (including, without limitation attorneys' and experts' fees and disbursements) of any kind whatsoever which may at any time be imposed upon, incurred by or asserted or awarded against client relating to, resulting from or arising out of (a) the breach of this agreement by this client, (b) the negligence of the client in handling, delivering or disclosing any hazardous substance, (c) the violation of the Client of any applicable law, (d) non-compliance by the Client with any environmental permit or (e) a material misrepresentation in disclosing the materials to be tested.

Hazard Disclosure.

Client represents and warrants that any sample delivered to LAB will be preceded or accompanied by complete written disclosure of the presence of any hazardous substances known or suspected by Client. Client further warrants that any sample containing any hazardous substance that is to be delivered to LAB will be packaged, labeled, transported, and delivered properly and in accordance with applicable laws.

Sample Handling.

Prior to LAB's acceptance of any sample (or after any revocation of acceptance), the entire risk of loss or of damage to such sample remains with Client. Samples are accepted when receipt is acknowledged on chain of custody documentation. In no event will LAB have any responsibility for the action or inaction of any carrier shipping or delivering any sample to or from LAB premises. Client authorizes LAB to proceed with the analysis of samples as received by the laboratory, recognizing that any samples not in compliance with all current DOH-ELAP-NELAP requirements for containers, preservation or holding time will be noted as such on the final report.

Disposal of hazardous waste samples is the responsibility of the Client. If the Client does not wish such samples returned, LAB may add storage and disposal fees to the final invoice. Maximum storage time for samples is 30 days after completion of analysis unless modified by applicable state or federal laws. Client will be required to give the LAB written instructions concerning disposal of these samples.

LAB reserves the absolute right, exercisable at any time, to refuse to receive delivery of, refuse to accept, or revoke acceptance of any sample, which, in the sole judgment of LAB (a) is of unsuitable volume, (b) may be or become unsuitable for or may pose a risk in handling, transport, or processing for any health, safety, environmental or other reason whether or not due to the presence in the sample of any hazardous substance, and whether or not such presence has been disclosed to LAB by Client or (c) if the condition or sample date make the sample unsuitable for analysis.

Legal Responsibility. LAB is solely responsible for performance of this contract, and no affiliated company, director, officer, employee, or agent shall have any legal responsibility hereunder, whether in contract or tort including negligence.

Assignment.

LAB may assign its performance obligations under this contract to other parties, as it deems necessary. LAB shall disclose to Client any assignee (subcontractor) by ELAP ID # on the submitted final report.

Force Majeure.

LAB shall have no responsibility or liability to the Client for any failure or delay in performance by LAB, which results in whole or in part from any cause or circumstance beyond the reasonable control of LAB. Such causes and circumstances shall include, but not limited to, acts of God, acts or orders of any government authority, strikes or other labor disputes, natural disasters, accidents, wars, civil disturbances, difficulties or delays in transportation, mail or delivery services, inability to obtain sufficient services or supplies from LAB's usual suppliers, or any other cause beyond LAB's reasonable control.

Law.

This contract shall be continued under the laws of the State of New York without regard to its conflicts of laws provision.

This report is part of a multipage document and should only be evaluated in its entirety. The Chain of Custody provides additional sample information, including compliance with the sample condition requirements upon receipt.

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		PARADIGM				PROJECT REFERENCE	i e	-	E E	соггестер	16:15	14:00	15145	(- 7	14:30	17.00	16:30		od Time	vility conting	X					ded:	
1		PA				RCF C				DATE COLLECTED	10/29/21	10/20/2	10/20/21	10/36/21	12/12/01	10/31/21	11/1/2	11/1/11	16/25/01	Turnaround Time	Availat	Standard 5 day	10 day	Rush 3 day	Rush 2 day	Rush 1 day	Date Needed please indicate date needed:	

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹								
Volatile Organics	Water	Soil							
Method 8260	(µg/L)	(µg/kg) ²							
Chloromethane	5	5							
Bromomethane	5	5							
Vinyl Chloride	5	5							
Chloroethane	5	5							
Methylene Chloride	3	3							
Acetone	5	5							
Carbon Disulfide	5	5							
1,1-Dichloroethylene	5	5							
1,1-Dichloroethane	5	5							
1,2-Dichloroethylene (total)	5	5							
Chloroform	5	5							
1,2-Dichloroethane	2	2							
2-Butanone	5	5							
1,1,1-Trichloroethane	5	5							
Carbon Tetrachloride	2	2							
Bromodichloromethane									
1,2-Dichloropropane	1	1							
	1	1							
cis-1,3-Dichloropropene	5	5							
Trichloroethane	5	5							
Dibromochloromethane	5	5							
1,1,2-Trichloroethane	3	3							
Benzene	1	11							
trans-1,3-Dichloropropene	5	5							
Bromoform	4	4							
4-Methyl-2-pentanone	5	5							
2-Hexanone	5	5							
Tetrachloroethene	1	1							
Toluene	5	5							
1,1,2,2-Tetrachloroethane	1	1							
Chlorobenzene	5	5							
Ethylbenzene	4	4							
Styrene	5	5							
2-Chloroethyl Vinyl Ether	5	5							
1,2-Dichlorobenzene	5	5							
1,3-Dichlorobenzene	5	5							
1,4-Dichlorobenzene	5	5							
Vinyl Acetate	5	5							
Total Xylenes	5	5							
Semivolatile Organics	Water	Soil							
Method 8270	(μg/L)	(µg/kg)							
1,2,4-Trichlorobenzene	1	33							
1,2-Dichlorobenzene	10	330							
1,2-Diphenylhydrazine	10	330							
1,3-Dichlorobenzene	10	330							
1,4-Dichlorobenzene	10	330							
1,4-Dioxane	10	330							
2,4,5-Trichlorophenol	10	330							
2,4,6-Trichlorophenol	10	330							
2,4-Dichlorophenol	10	330							

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Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹				
Semivolatile Organics	Water	Soil			
Method 8270 (Cont'd.)	(µg/L)	(µg/kg)			
2,4-Dimethylphenol	10	330			
2,4-Dinitrophenol	40	1300			
2,4-Dinitrotoluene	2	67			
2,6-Dinitrotoluene	2	67			
2-Chloronaphthalene	10	330			
2-Chlorophenol	10	330			
2-Methylnaphthalene	10	330			
2-Methylphenol	10	330			
2-Nitroaniline	20	670			
2-Nitrophenol	10	330			
3,3'-Dichlorobenzidene	20	670			
3-Nitroaniline	20	670			
4,6-Dinitro-2-methylphenol	40	1300			
4-Bromophenyl-phenylether	10	330			
4-Chloro-3-methylphenol	10	330			
4-Chloroaniline	10	330			
4-Chlorophenyl-phenylether	10	330			
4-Methylphenol	10	330			
4-Nitroaniline	20	670			
4-Nitrophenol	40	1300			
Acenaphthene	10	330			
Acenaphthylene	10	330			
Acetophenone	10	330			
Aniline	10	330			
Anthracene	10	330			
Atrazine	10	330			
Benzaldehyde	10	330			
Benzidine	40	1300			
Benzo(a)anthracene	1	33			
Benzo(a)pyrene	1	33			
Benzo(b)fluoranthene	1	33			
Benzo(g,h,i)perylene	10	330			
Benzo(k)fluoranthene	1	33			
Benzoic Acid	10	330			
Benzyl Alcohol	10	330			
bis(2-chloroethoxy)methane	10	330			
bis(2-chloroethyl)ether	1	33			
bis(2-chloroisopropyl)ether	10	330			
bis(2-ethylhexyl)phthalate	10	330			
Butylbenzylphthalate	10	330			
Caprolactam	10	330			
Carbazole	10	330			
Chrysene	10	330			
Dibenzo(a,h)anthracene	1	33			
Dibenzofuran	10	330			
Diethylphthalate	10	330			
Dimethylphthalate	10	330			
Di-n-butyl phthalate	10	330			
Di-n-octyl phthalate	10				
Difficulty philialate	10	330			

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Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹				
Semivolatiles	Water	Soil			
Method 8270 (Cont'd.)	(µg/L)	(µg/kg)			
Diphenyl	10	330			
Fluoranthene	10	330			
Fluorene	10	330			
Hexachlorobenzene	1	33			
Hexachlorobutadiene	2	67			
Hexachlorocyclopentadiene	10	330			
Hexachloroethane	1	33			
Indeno(1,2,3-cd)pyrene	1	33			
Isophorone	10	330			
N,N-Dimethylaniline	1	33			
Naphthalene	10	330			
Nitrobenzene	1	33			
N-Nitrosodimethylamine	10	330			
N-Nitroso-di-n-propylamine	1	33			
N-Nitrosodiphenylamine	10	330			
Pentachlorophenol	40	1300			
Phenanthrene	10	330			
Phenol	10	330			
Pyrene	10	330			
Pyridine	10	330			
TAL Metals (6010/7470)	Water	Soil			
Aluminum	(µg/L)	(µg/kg)			
		40			
Antimony	N-4	2			
Arsenic		1			
Barium		40			
Beryllium		0.4			
Cadmium		1 1			
Calcium	A.P.	1000			
Chromium		2			
Cobalt	E-	10			
Copper	**	5			
ron		30			
Lead		1			
Magnesium	×=	1000			
Manganese		3			
Mercury		0.033			
Nickel		8			
Potassium		1000			
Selenium		1			
Silver		2			
Sodium		1000			
Thallium		2			
Vanadium		10			
Zinc		6			
Supplemental Parameters	Water	Soil			
	(µg/L)	(mg/kg)			
Total Organic Carbon (Lloyd Kahn)		(mg/kg) 100			

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Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹			
Supplemental Parameters (Cont'd.)	Water (µg/L)	Soil (mg/kg)		
Nitrate Method 353.2	100			
Ammonia Method 350.1	100			
Iron Method 200.7	150			
Manganese Method 200.7	15			
Sulfate Method 375.4	5,000			
Sulfide Method 376.1	1,000			
Orthophosphate Method 365.2	30	-		
Alkalinity Method 310.1	5,000			
Methane Method 3810				
Reactive Sulfide		20		
Reactive Cyanide		25		
TCLP Benzene		1		
Total Sulfur		50		
Chemical Oxygen Demand		120		

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Notes:

μg/L = micrograms per liter

μg/kg = micrograms per kilogram

mg/kg = milligrams per kilogram

¹ Specific quantitation limits are highly matrix dependent. The quantitation limits listed are for guidance and may not always be achievable due to matrix interference.

² Quantitation limits for source materials/soil/sediment are based on wet weight. The quantitation limits calculated by the laboratory for source materials/soil/sediment (calculated on a dry weight basis) will be higher.



Chain of Custody Supplement

Client:	Neu-Velle	Completed by:	MolyVail
Lab Project ID:	214958	Date:	1/12/2/
	Sample Cond Per NELAC/ELA	ition Requirements P 210/241/242/243/244	
Condition	NELAC compliance with the sam Yes	ple condition requirements upon No	receipt- N/A
Container Type			
Comments			- N
Transferred to method- compliant container			
Headspace (<1 mL) Comments	VOA		SVOA
Preservation Comments	NOA		500
Chlorine Absent (<0.10 ppm per test strip) Comments			
Holding Time Comments			
Temperature Comments	y°cial		
Compliant Sample Quantity/Ty	ре		
Comments			



ANALYTICAL REPORT

Lab Number: L2160069

Client: Paradigm Environmental Services

179 Lake Avenue Rochester, NY 14608

ATTN: Jane Daloia Phone: (585) 647-2530

Project Name: 214958
Project Number: 214958
Report Date: 11/09/21

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2160069-01	GEN-MW4-102921 214958-01	WATER	Not Specified	10/29/21 16:15	11/02/21
L2160069-02	GEN-MW1-103021 214958-02	WATER	Not Specified	10/30/21 14:00	11/02/21
L2160069-03	GEN-MW2-103021 214958-03	WATER	Not Specified	10/30/21 15:45	11/02/21
L2160069-04	GEN-DUP-103021 214958-04	WATER	Not Specified	10/30/21 00:00	11/02/21
L2160069-05	GEN-MW3-103121 214958-05	WATER	Not Specified	10/31/21 12:30	11/02/21
L2160069-06	GEN-MW6-103121 214958-06	WATER	Not Specified	10/31/21 14:30	11/02/21
L2160069-07	GEN-EB-110121 214958-07	WATER	Not Specified	11/01/21 12:00	11/02/21
L2160069-08	GEN-MW8-110121 214958-08	WATER	Not Specified	11/01/21 15:00	11/02/21
L2160069-09	GEN-MW7-110121 214958-09	WATER	Not Specified	11/01/21 16:30	11/02/21



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Cattlin Wallet Caitlin Walukevich

Authorized Signature:

Title: Technical Director/Representative

ΔLPHA

Date: 11/09/21

ORGANICS



SEMIVOLATILES



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-01 Date Collected: 10/29/21 16:15

Client ID: GEN-MW4-102921 214958-01 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/04/21 20:06

Analyst: ALS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM -	Semivolatile Organics by GC/MS-SIM - Westborough Lab						
Acenaphthene	ND		ug/l	0.10	0.04	1	
Fluoranthene	ND		ug/l	0.10	0.04	1	
Naphthalene	ND		ug/l	0.10	0.04	1	
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1	
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1	
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1	
Chrysene	ND		ug/l	0.10	0.04	1	
Acenaphthylene	ND		ug/l	0.10	0.04	1	
Anthracene	ND		ug/l	0.10	0.04	1	
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1	
Fluorene	ND		ug/l	0.10	0.04	1	
Phenanthrene	ND		ug/l	0.10	0.02	1	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1	
Pyrene	ND		ug/l	0.10	0.04	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	70	23-120	
2-Fluorobiphenyl	69	15-120	
4-Terphenyl-d14	67	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-02 Date Collected: 10/30/21 14:00

Client ID: GEN-MW1-103021 214958-02 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 17:19

Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM - Westborough Lab						
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.02	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.02	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.02	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
ND		ug/l	0.10	0.04	1	
	ND N	ND N	ND ug/l ND ug/l	ND	ND	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	72	23-120	
2-Fluorobiphenyl	75	15-120	
4-Terphenyl-d14	79	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-03 Date Collected: 10/30/21 15:45

Client ID: GEN-MW2-103021 214958-03 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 17:35

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM - Westborough Lab						
Acenaphthene	ND		ug/l	0.10	0.04	1
Fluoranthene	ND		ug/l	0.10	0.04	1
Naphthalene	ND		ug/l	0.10	0.04	1
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1
Chrysene	ND		ug/l	0.10	0.04	1
Acenaphthylene	ND		ug/l	0.10	0.04	1
Anthracene	ND		ug/l	0.10	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1
Fluorene	ND		ug/l	0.10	0.04	1
Phenanthrene	0.02	J	ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1
Pyrene	ND		ug/l	0.10	0.04	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	77	23-120	
2-Fluorobiphenyl	77	15-120	
4-Terphenyl-d14	79	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-04 Date Collected: 10/30/21 00:00

Client ID: GEN-DUP-103021 214958-04 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 17:52

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		/1	0.10	0.04	1	
·			ug/l				
Fluoranthene	ND		ug/l	0.10	0.04	1	
Naphthalene	ND		ug/l	0.10	0.04	1	
Benzo(a)anthracene	0.02	J	ug/l	0.10	0.02	1	
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1	
Benzo(b)fluoranthene	0.03	J	ug/l	0.10	0.02	1	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1	
Chrysene	ND		ug/l	0.10	0.04	1	
Acenaphthylene	ND		ug/l	0.10	0.04	1	
Anthracene	ND		ug/l	0.10	0.04	1	
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1	
Fluorene	ND		ug/l	0.10	0.04	1	
Phenanthrene	0.02	J	ug/l	0.10	0.02	1	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1	
Pyrene	ND		ug/l	0.10	0.04	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	66	23-120	
2-Fluorobiphenyl	74	15-120	
4-Terphenyl-d14	82	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-05 Date Collected: 10/31/21 12:30

Client ID: GEN-MW3-103121 214958-05 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 18:08

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM	- Westborough Lal	b				
Acenaphthene	ND		ug/l	0.10	0.04	1
Fluoranthene	ND		ug/l	0.10	0.04	1
Naphthalene	ND		ug/l	0.10	0.04	1
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1
Chrysene	ND		ug/l	0.10	0.04	1
Acenaphthylene	ND		ug/l	0.10	0.04	1
Anthracene	ND		ug/l	0.10	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1
Fluorene	ND		ug/l	0.10	0.04	1
Phenanthrene	ND		ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1
Pyrene	ND		ug/l	0.10	0.04	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	58	23-120	
2-Fluorobiphenyl	63	15-120	
4-Terphenyl-d14	77	41-149	



Extraction Method: EPA 3510C

11/03/21 15:38

Extraction Date:

Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-06 D Date Collected: 10/31/21 14:30

Client ID: GEN-MW6-103121 214958-06 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8270D-SIM
Analytical Date: 11/09/21 11:40

Analyst: RP

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM	Westborough La	ab				
Acenaphthene	1.7		ug/l	1.0	0.35	10
Fluoranthene	ND		ug/l	1.0	0.38	10
Naphthalene	170		ug/l	1.0	0.43	10
Benzo(a)anthracene	ND		ug/l	1.0	0.18	10
Benzo(a)pyrene	ND		ug/l	1.0	0.39	10
Benzo(b)fluoranthene	ND		ug/l	1.0	0.16	10
Benzo(k)fluoranthene	ND		ug/l	1.0	0.42	10
Chrysene	ND		ug/l	1.0	0.38	10
Acenaphthylene	18		ug/l	1.0	0.35	10
Anthracene	0.46	J	ug/l	1.0	0.35	10
Benzo(ghi)perylene	ND		ug/l	1.0	0.42	10
Fluorene	4.0		ug/l	1.0	0.37	10
Phenanthrene	1.6		ug/l	1.0	0.15	10
Dibenzo(a,h)anthracene	ND		ug/l	1.0	0.39	10
Indeno(1,2,3-cd)pyrene	ND		ug/l	1.0	0.40	10
Pyrene	ND		ug/l	1.0	0.40	10

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	75	23-120	
2-Fluorobiphenyl	86	15-120	
4-Terphenyl-d14	107	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-07 Date Collected: 11/01/21 12:00

Client ID: GEN-EB-110121 214958-07 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 18:41

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM	- Westborough La	ab				
Acenaphthene	ND		ug/l	0.10	0.04	1
Fluoranthene	ND		ug/l	0.10	0.04	1
Naphthalene	ND		ug/l	0.10	0.04	1
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1
Chrysene	ND		ug/l	0.10	0.04	1
Acenaphthylene	ND		ug/l	0.10	0.04	1
Anthracene	ND		ug/l	0.10	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1
Fluorene	ND		ug/l	0.10	0.04	1
Phenanthrene	ND		ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1
Pyrene	ND		ug/l	0.10	0.04	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	85	23-120	
2-Fluorobiphenyl	84	15-120	
4-Terphenyl-d14	83	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-08 Date Collected: 11/01/21 15:00

Client ID: GEN-MW8-110121 214958-08 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 18:58

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Aggraphthana	ND		/1	0.10	0.04	1	
Acenaphthene			ug/l			I	
Fluoranthene	ND		ug/l	0.10	0.04	1	
Naphthalene	0.22		ug/l	0.10	0.04	1	
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1	
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1	
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1	
Chrysene	ND		ug/l	0.10	0.04	1	
Acenaphthylene	ND		ug/l	0.10	0.04	1	
Anthracene	ND		ug/l	0.10	0.04	1	
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1	
Fluorene	ND		ug/l	0.10	0.04	1	
Phenanthrene	0.04	J	ug/l	0.10	0.02	1	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1	
Pyrene	ND		ug/l	0.10	0.04	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	67	23-120	
2-Fluorobiphenyl	69	15-120	
4-Terphenyl-d14	77	41-149	



Project Name: 214958 Lab Number: L2160069

Project Number: 214958 Report Date: 11/09/21

SAMPLE RESULTS

Lab ID: L2160069-09 Date Collected: 11/01/21 16:30

Client ID: GEN-MW7-110121 214958-09 Date Received: 11/02/21 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 11/03/21 15:38
Analytical Date: 11/05/21 19:14

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM -	Westborough La	ab				
Acenaphthene	ND		ug/l	0.10	0.04	1
Fluoranthene	ND		ug/l	0.10	0.04	1
Naphthalene	ND		ug/l	0.10	0.04	1
Benzo(a)anthracene	0.02	J	ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1
Benzo(b)fluoranthene	0.02	J	ug/l	0.10	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1
Chrysene	ND		ug/l	0.10	0.04	1
Acenaphthylene	ND		ug/l	0.10	0.04	1
Anthracene	ND		ug/l	0.10	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1
Fluorene	ND		ug/l	0.10	0.04	1
Phenanthrene	0.02	J	ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1
Pyrene	ND		ug/l	0.10	0.04	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	44	23-120	
2-Fluorobiphenyl	50	15-120	
4-Terphenyl-d14	63	41-149	



 Project Name:
 214958

 Lab Number:
 L2160069

Project Number: 214958 Report Date: 11/09/21

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D-SIM Extraction Method: EPA 3510C
Analytical Date: 11/04/21 19:16 Extraction Date: 11/03/21 15:38

Analyst: ALS

Acenaphthene ND ug/l 0.10 0.04 Fluoranthene ND ug/l 0.10 0.04 Fluoranthene ND ug/l 0.10 0.04 Naphthalene ND ug/l 0.10 0.04 Benzo(a)anthracene ND ug/l 0.10 0.02 Benzo(a)pyrene ND ug/l 0.10 0.04 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04 <th>arameter</th> <th>Result</th> <th>Qualifier Units</th> <th>RL</th> <th>MDL</th> <th></th>	arameter	Result	Qualifier Units	RL	MDL	
Fluoranthene ND ug/l 0.10 0.04 Naphthalene ND ug/l 0.10 0.04 Benzo(a)anthracene ND ug/l 0.10 0.02 Benzo(a)pyrene ND ug/l 0.10 0.04 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	emivolatile Organics by GC	C/MS-SIM - Westbo	rough Lab for samp	le(s): 01-09	Batch: Wo	G1566746-1
Naphthalene ND ug/l 0.10 0.04 Benzo(a)anthracene ND ug/l 0.10 0.02 Benzo(a)pyrene ND ug/l 0.10 0.04 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Acenaphthene	ND	ug/l	0.10	0.04	
Benzo(a)anthracene ND ug/l 0.10 0.02 Benzo(a)pyrene ND ug/l 0.10 0.04 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Fluoranthene	ND	ug/l	0.10	0.04	
Benzo(a)pyrene ND ug/l 0.10 0.04 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Naphthalene	ND	ug/l	0.10	0.04	
Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Benzo(a)anthracene	ND	ug/l	0.10	0.02	
Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Benzo(a)pyrene	ND	ug/l	0.10	0.04	
Chrysene ND ug/l 0.10 0.04 Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Benzo(b)fluoranthene	ND	ug/l	0.10	0.02	
Acenaphthylene ND ug/l 0.10 0.04 Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Benzo(k)fluoranthene	ND	ug/l	0.10	0.04	
Anthracene ND ug/l 0.10 0.04 Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Chrysene	ND	ug/l	0.10	0.04	
Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Acenaphthylene	ND	ug/l	0.10	0.04	
Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Anthracene	ND	ug/l	0.10	0.04	
Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Benzo(ghi)perylene	ND	ug/l	0.10	0.04	
Dibenzo(a,h)anthracene ND ug/l 0.10 0.04 Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Fluorene	ND	ug/l	0.10	0.04	
Indeno(1,2,3-cd)pyrene ND ug/l 0.10 0.04	Phenanthrene	ND	ug/l	0.10	0.02	
	Dibenzo(a,h)anthracene	ND	ug/l	0.10	0.04	
Pyrene ND ug/l 0.10 0.04	Indeno(1,2,3-cd)pyrene	ND	ug/l	0.10	0.04	
	Pyrene	ND	ug/l	0.10	0.04	

		Acceptance
Surrogate	%Recovery	Qualifier Criteria
O Florent cod	40	04.400
2-Fluorophenol	43	21-120
Phenol-d6	29	10-120
Nitrobenzene-d5	72	23-120
2-Fluorobiphenyl	71	15-120
2,4,6-Tribromophenol	70	10-120
4-Terphenyl-d14	67	41-149



Lab Control Sample Analysis Batch Quality Control

Project Name: 214958
Project Number: 214958

Lab Number: L2160069

Report Date:

11/09/21

	LCS		LCSD		%Recove	ry		RPD	
arameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual	Limits	
emivolatile Organics by GC/MS-SIM -	Westborough Lab A	Associated samp	ole(s): 01-09	Batch:	WG1566746-2	WG1566746-3			
Acenaphthene	72		69		40-140	4		40	
Fluoranthene	78		69		40-140	12		40	
Naphthalene	74		69		40-140	7		40	
Benzo(a)anthracene	80		70		40-140	13		40	
Benzo(a)pyrene	81		70		40-140	15		40	
Benzo(b)fluoranthene	78		70		40-140	11		40	
Benzo(k)fluoranthene	80		67		40-140	18		40	
Chrysene	78		69		40-140	12		40	
Acenaphthylene	84		79		40-140	6		40	
Anthracene	78		71		40-140	9		40	
Benzo(ghi)perylene	80		71		40-140	12		40	
Fluorene	73		70		40-140	4		40	
Phenanthrene	75		68		40-140	10		40	
Dibenzo(a,h)anthracene	82		74		40-140	10		40	
Indeno(1,2,3-cd)pyrene	80		70		40-140	13		40	
Pyrene	79		70		40-140	12		40	



L2160069

Lab Control Sample Analysis

Project Name: 214958

214958

Project Number:

Batch Quality Control

_ . _

Report Date: 11/09/21

Lab Number:

LCS LCSD %Recovery RPD Parameter %Recovery Qual %Recovery Qual Limits RPD Qual Limits

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-09 Batch: WG1566746-2 WG1566746-3

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
2-Fluorophenol	48	46	21-120
Phenol-d6	33	32	10-120
Nitrobenzene-d5	80	75	23-120
2-Fluorobiphenyl	76	72	15-120
2,4,6-Tribromophenol	118	76	10-120
4-Terphenyl-d14	64	67	41-149



Matrix Spike Analysis Batch Quality Control

Project Name: 214958 **Project Number:** 214958 Lab Number:

L2160069

Report Date:

11/09/21

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery	Recover Qual Limits	y RPD	RPD Qual Limits
Semivolatile Organics by Client ID: GEN-MW4-102	GC/MS-SIM - We	stborough Lab	Associated	d sample(s): 01-	09 QC	Batch ID:	WG1566746-4	WG1566746-5		ple: L2160069-01
Acenaphthene	ND	10	6.8	68		6.9	69	40-140	1	40
Fluoranthene	ND	10	6.8	68		7.2	72	40-140	6	40
Naphthalene	ND	10	6.8	68		7.1	71	40-140	4	40
Benzo(a)anthracene	ND	10	7.1	71		7.4	74	40-140	4	40
Benzo(a)pyrene	ND	10	6.9	69		7.3	73	40-140	6	40
Benzo(b)fluoranthene	ND	10	7.0	70		7.2	72	40-140	3	40
Benzo(k)fluoranthene	ND	10	6.9	69		7.2	72	40-140	4	40
Chrysene	ND	10	6.7	67		6.9	69	40-140	3	40
Acenaphthylene	ND	10	7.8	78		7.9	79	40-140	1	40
Anthracene	ND	10	6.9	69		7.1	71	40-140	3	40
Benzo(ghi)perylene	ND	10	7.1	71		7.5	75	40-140	5	40
Fluorene	ND	10	6.9	69		7.0	70	40-140	1	40
Phenanthrene	ND	10	6.7	67		6.8	68	40-140	1	40
Dibenzo(a,h)anthracene	ND	10	7.4	74		7.8	78	40-140	5	40
ndeno(1,2,3-cd)pyrene	ND	10	7.1	71		7.4	74	40-140	4	40
Pyrene	ND	10	6.7	67		7.3	73	40-140	9	40

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
2,4,6-Tribromophenol	74	78	10-120
2-Fluorobiphenyl	72	75	15-120
2-Fluorophenol	46	53	21-120
4-Terphenyl-d14	65	74	41-149



Matrix Spike Analysis Batch Quality Control

Project Name: 214958
Project Number: 214958

Lab Number:

L2160069

Report Date:

11/09/21

	Native	MS	MS	MS		MSD	MSD		Recovery			RPD
Parameter	Sample	Added	Found	%Recovery	Qual	Found	%Recovery	Qual	Limits	RPD	Qual	Limits

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-09 QC Batch ID: WG1566746-4 WG1566746-5 QC Sample: L2160069-01 Client ID: GEN-MW4-102921 214958-01

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
- Nitrobenzene-d5	75	79	23-120
Phenol-d6	33	38	10-120



Lab Number: L2160069

Report Date: 11/09/21

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Project Name:

Project Number: 214958

Cooler Custody Seal

214958

A Absent

Container Info	rmation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	pН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2160069-01A	Amber 1000ml unpreserved	Α	8	8	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-01A1	Amber 1000ml unpreserved	Α	9	9	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-01A2	Amber 1000ml unpreserved	Α	9	9	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-02B	Amber 1000ml unpreserved	Α	7	7	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-03B	Amber 1000ml unpreserved	Α	8	8	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-04B	Amber 1000ml unpreserved	Α	8	8	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-05B	Amber 1000ml unpreserved	Α	8	8	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-06B	Amber 1000ml unpreserved	Α	9	9	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-07B	Amber 1000ml unpreserved	Α	9	9	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-08B	Amber 1000ml unpreserved	Α	8	8	2.7	Υ	Absent		NYTCL-8270-SIM(7)
L2160069-09B	Amber 1000ml unpreserved	Α	9	9	2.7	Υ	Absent		NYTCL-8270-SIM(7)



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

GLOSSARY

Acronyms

EDL

LOQ

MS

RL

RPD

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).

EMPC - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

 Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

 Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

 NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

NR - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.

Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEQ - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



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 Lab Number:
 L2160069

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 Report Date:
 11/09/21

Footnotes

 The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

1

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benza(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

- A -Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



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 Report Date:
 11/09/21

Data Qualifiers

- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q -The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
 (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:17873 Revision 19

Page 1 of 1

Published Date: 4/2/2021 1:14:23 PM

Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

EPA 8260C/8260D: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522, EPA 537.1.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Pre-Qualtrax Document ID: 08-113 Document Type: Form

179 Lake Avenue, Rochester, NY 14608 Office (585) 647-2530 Fax (585) 647-3311

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DATA USABILITY SUMMARY REPORT (DUSR)

RGE Geneseo Former MGP Site

SDGs: 214958

8 Water Samples, equipment blank and trip blank

Prepared for:

Neu Velle, LLC 1667 Lake Ave., Bldg. 59, Suite 101 Rochester, NY 14615 **Attention: Kyle Miller**

February 2022



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Table 4-1 Data Validation Guidance Documents

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Summaries of Validated Results

Table 6-1 8260-BTEX Table 6-2 8270-SIM-PAHs

REVIEWER'S NARRATIVE

Neu-Velle SDG 214958: RGE Geneseo Former MGP Site

The data associated with this Sample Delivery Group (SDG) 214958, analyzed by Paradigm Environmental Services, Inc. Rochester, NY have been reviewed in accordance with assessment criteria provided by the New York State Department of Environmental Conservation following the review procedures provided in the USEPA Functional Guidelines for evaluating organic and inorganic data.

All analytical results reported by the laboratory are considered valid and acceptable except results that have been qualified as rejected, "R". Results qualified as estimated "J", or as non-detects, "U", are considered usable for the purpose of evaluating water and/or soil quality. However, these qualifiers indicate that the accuracy and/or precision of the analytical result is questionable. A summary of all data that have been qualified and the reasons for qualification are provided in the following data usability summary report (DUSR).

Two facts should be noted by all data users. First, the "R" qualifier means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the analyte is present or not. Values qualified with an "R" should not appear on the final data tables because they cannot be relied upon, even as the last resort. Second, no analyte concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error.

Reviewer's Signature:	Míchael K. Perry	Date:	2/16/2022	
	Michael K. Perry			
	Chemist			

1.0 SUMMARY

SITE: RGE Geneseo

Former MGP Site

SAMPLING DATE: October 29 – November 01, 2021

SAMPLE TYPE: 8 water samples, equipment and trip blank

LABORATORY: Paradigm Environmental Services, Inc.

Rochester, NY

SDG No.: 214958

2.0 INTRODUCTION

This data usability summary report (DUSR) was prepared in accordance with guidance provided by the New York State Department of Environmental Conservation (NYSDEC). The DUSR is based on a review and evaluation of the laboratory analytical data package. Specifically, the NYSDEC guidance recommends review and evaluation of the following elements of the data package:

- Completeness of the data package as defined under the requirements of the NYSDEC Analytical Services Protocols (ASP) Category B or the United States Environmental Protection Agency (USEPA) Contract Laboratory Program (CLP) deliverables,
- Compliance with established analyte holding times,
- Adherence to quality control (QC) limits and specifications for blanks, instrument tuning and calibration, surrogate recoveries, spike recoveries, laboratory duplicate analyses, and other QC criteria,
- Adherence to established analytical protocols,
- Conformance of data summary sheets with raw analytical data, and
- Use of correct data qualifiers.

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Data deficiencies, analytical protocol deviations, and quality control problems identified using the review criteria above and their effect on the analytical results are discussed in this report.

3.0 SAMPLE AND ANALYSIS SUMMARY

The data packages consists of analytical results for eight water samples, equipment blank, and a trip blank collected on October 29 – November 01, 2021. These samples were analyzed for volatile organic compounds (BTEX) and Semi-Volatile Organic Compounds (SVOCs).

All analyses were performed by Paradigm Environmental Services, Inc., Rochester, NY and analyzed as SDG: 214958. The SVOCs were subcontracted to Alpha Analytical in Westborough, MA and analyzed as SDG: L2160069 for PAHs by 8270-D-SIM. The analytical results were provided in NYSDEC ASP Category B format, which includes all raw analytical data and laboratory QC data.

4.0 GUIDANCE DOCUMENTS AND DATA REVIEW CRITERIA

The guidance documents appropriate for reviewing laboratory quality control (QC) data and assigning data qualifiers (flags) to analytical results were selected from those listed in Table 4-1. The QC limits established in the documents applicable to this data review were used to assess the quality of the analytical results. In some cases, however, QC limits established internally by the laboratory were taken into account to determine data quality.

The QC criteria considered for assessing the usability of the reported analytical results provided for each analyte type (i.e. VOCs, SVOCs, metals, etc.) are listed in Table 4-2. These criteria may vary with the analytical method utilized by the laboratory. These criteria comply with the guidance recommended in Section 2.0 above.

5.0 DATA VALIDATION QUALIFIERS

The letter qualifiers (flags) used to define data usability are described briefly below. These letters are assigned by the data validator to analytical results having questionable accuracy and/or precision as determined by reviewing the laboratory QC data associated with the analytical results.

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TABLE 4-1

Guidance Used For Validating Laboratory Analytical Data

Analyte Group	Guidance	Date
Metals (ICP-AES)	USEPA SOP HW-3a, Rev. 1	September 2016
Metals (Hg & CN)	USEPA SOP HW-3c, Rev. 1	September 2016
Volatile Organic Compounds (by Methods 8260B & 8260C)	USEPA SOP HW-24, Rev. 4	September 2014
Semi-Volatile Organic Compounds (by Method 8270D)	USEPA SOP HW-22 Rev. 5	December 2010
Pesticides (by Method 8181B)	USEPA SOP HW-44, Rev. 1.1	December 2010
Chlorinated Herbicides (by Method 8151A)	USEPA SOP HW-17, Rev. 3.1	December 2010
Polychlorinated Biphenyls (PCBs)	USEPA SOP HW-37A, Rev. 0	June 2015
Volatile Organic Compounds (Air) (by Method TO-15)	USEPA SOP HW-31, Rev. 6	September 2016
Per- and PolyFluoroAlkyl Substances (PFAS)	* NYSDEC	January 2021
General Chemistry Parameters	per NYSDEC ASP	July 2005

^{*} Sampling, Analysis, and Assessment of Per- and Polyfluoroalkyl Substances (PFAS) Under NYSDEC's Part 375 Remedial Programs, Appendix I

TABLE 4-2

QUALITY CONTROL CRITERIA USED FOR VALIDATING LABORATORY ANALYTICAL DATA

VOCs	SVOCs	Pesticides/PCBs	Metals	Gen Chemistry	PFAS
Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg
Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation
Holding Time	Holding Time	Holding Time	Holding Time	Holding Times	Holding Time
System Monitoring	Surrogate Recoveries	Surrogate Recoveries	Initial/Continuing	Calibration	Instr Performance
Compounds	Lab Control Sample	Matrix Spikes	Calibration	Lab Control Samples	Check
Lab Control Sample	Matrix Spikes	Blanks	CRDL Standards	Blanks	Initial Calibration
Matrix Spikes	Blanks	Instrument Calibration	Blanks	Spike Recoveries	Continuing Calibration
Blanks	Instrument Tuning	& Verification	Interference Check	Lab Duplicates	Blanks
Instrument Tuning	Internal Standards	Comparison of	Sample		Surrogates
Internal Standards	Initial Calibration	duplicate	Spike Recoveries		Lab Fortified Blank
Initial Calibration	Continuing Calibration	GC column results	Lab Duplicate		Matrix Spikes
Continuing Calibration	Lab Qualifiers	Analyte ID	Lab Control Sample		Internal Standards
Lab Qualifiers	Field Duplicate	Lab Qualifiers	ICP Serial Dilutions		
Field Duplicate		Field Duplicate	Lab Qualifiers		
			Field Duplicate		

Method TO-15 (Air)

Completeness of Pkg
Sample Preservation
Holding Time
Canister Certification
Instrument Tuning
Initial Calibration and
Instrument Performance
Daily Calibration
Blanks
Lab Control Sample
Field Duplicate

The laboratory may also use various letters and symbols to flag analytical results generated when QC limits were exceeded. The meanings of these flags may differ from those used by the independent data validator. Those used by the laboratory are provided with the analytical results.

NOTE: The assignment of data qualifiers by the data reviewer (validator) to laboratory analytical results should not necessarily be interpreted by the data user as a measure of laboratory ability or proficiency. Rather, the qualifiers are intended to provide a measure of data accuracy and precision to the data user, which, for example, may provide a level of confidence in determining whether or not standards or cleanup objectives have been met.

- U The analyte was analyzed for but was not detected at or above the sample quantitation limit.
- J The analyte was positively identified; the associated numerical value is the *approximate* concentration of the analyte in the sample. (The magnitude of any \pm value associated with the result is not determined by data validation).
- **J**+ The result is an estimated quantity and may be biased high.
- **J-** The result is an estimated quantity and may be biased low.
- **UJ** The analyte was analyzed for but not detected. The reported quantitation limit is approximate and may inaccurate or imprecise.
- R The sample result is rejected (i.e., is unusable) due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- **NJ** The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

The validated analytical results are attached to this report. Validation qualifiers (flags) are indicated in red print. Data sheets having qualified data are signed and dated by the data reviewer.

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6.0 RESULTS OF THE DATA REVIEW

The results of the data review are summarized in Tables 6-1 through 6-2. The tables list the samples where QC criteria were found to exceed acceptable limits and the actions taken to qualify the associated analytical results.

7.0 TOTAL USABLE DATA

For SDG 214958, eight samples, equipment blank and a trip blank were analyzed and results were reported for 204 analytes. Even though some results were flagged with a "J" as estimated, all results (100%) are considered usable.

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SDG 214958

Table 6-1 8260 - BTEX

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
GEN-MW4 GEN-EB GEN-MW8 Trip Blank	All analytes	J detects	PFB > QC limit	Detected data are estimated
GEN-MW4	Benzene Ethyl benzene Toluene	J detects	MS/MSD > QC limit	Detected data are estimated
GEN-MW8	Benzene	J detect	Detected in equipment blank	Detected data are estimated

Table 6-2 8270-SIM-PAH

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
none		none		

ACRONYMS

BSP Blank Spike

CCAL Continuing Calibration

CCB Continuing Calibration Blank

CCV Continuing Calibration Verification

CRDL Contract Required Detection Limit

CRQL Contract Required Quantitation Limit

% D Percent Difference

ICAL Initial Calibration

ICB Initial Calibration Blank

IS Internal Standard

LCS Laboratory Control Sample

MS/MSD Matrix Spike/Matrix Spike Duplicate

QA Quality Assurance

QC Quality Control

%R Percent recovery

RPD Relative Percent Difference

RRF Relative Response Factor

% RSD Percent Relative Standard Deviation

TAL Target Analyte List (metals)

TCL Target Compound List (organics)

Appendix A

Validated Analytical Results

LAB PROJECT NARRATIVE: 214958 PROJECT NAME: RGE Geneseo Fmr. MGP Site

SDG: 4958-01 CLIENT: Neu-Velle

Seven groundwater samples were collected by the client between October 29 and November 1, 2021 and were received by the Paradigm laboratory on November 2, 2021. Samples were accompanied by a field duplicate, equipment blank, and trip blank. The samples were received under the conditions as noted on the chain-of-custody supplement. The samples were submitted with the Chains-of-Custody requesting the BTEX list for Volatiles and PAH by 8270SIM. All analyses, where applicable, were performed using EPA SW-846 Methods and the associated holding times.

The items noted in this case narrative address compliance with the referenced methods, NYSDOH ELAP rules, and any project specific data quality requirements. These may be different from the usability criteria referenced in any "Functional Guidelines" or other data review standards used by data validators.

GENERAL NOTES

ALL ANALYSES

The initial and continuing calibration reports are only evaluated for compounds that are on the sample summary report.

Regarding results on QC summary forms versus included raw data, due to calculations made at the instrument where many significant figures may be used, there may be slight discrepancies between the summary report result and that recorded on the raw data. This does not affect data usability.

VOLATILES

Regarding initial calibrations, it should be noted that the Quantitation Report concentrations supplied for the initial calibration reflect the calibration prior to updating. The response factors and areas are correct.

Regarding Quantitation Reports, it should be noted that the "#" symbol that appears on some of the Quantitation Reports is a software artifact and should be disregarded.

Compounds flagged with an "*" on the summary table have been calibrated using a non-average Response Factor calibration curve. The supporting curves are located after the initial calibration table.

Holding times were met for the samples.

Some of the surrogate recoveries for the samples and associated QC were not within acceptance limits (recoveries were high). Any outliers have been flagged with an "*" on the summary form and the sample reports.

Site specific QC was requested on GEN-MW4-102921 and there were outliers. These outliers have been flagged with an "*" on the summary form and an "M" on the sample reports. Matrix interference is suspected. The Laboratory Control Samples recovered within acceptance limits.

The Method Blanks were free from contamination within reportable ranges.

The instrument tunes passed all criteria and samples were within a 12-hour window.

The internal standards areas and retention times were within acceptance ranges for the samples and QC.

All data for the initial calibration was within acceptance limits for the reported analytes.

All continuing calibration data was within acceptance limits for the reported analytes.

Subcontracted Analyses

PAHs by EPA 8270D-SIM was sent to Alpha Analytical of Westborough, MA. Their reports are provided in their entirety as a separate entity after the Paradigm Environmental Services, Inc. report. Separate case narratives addressing the above parameters are included with their reports.

(signed) Steven DeVito Steven DeVito – Technical Director

BATCH LOG

Lab Name: Paradigm Environmental Services

214958 Neu-Velle Lab Project #: Client Name:

RGE Geneseo Fmr. MGP Site
N/A Client Project Name:

Client Project #: <u>4958-01</u> SDG No.:

Batch Due Date: 12/2/2021 Protocol: SW846 Report Due Date: 11/17/2021

LAB	MATRIX	CLIENT	REQUESTED ANALYSIS	DATE	DATE
SAMPLE NO.		SAMPLE ID		SAMPLED	REC'D
214958-01	Groundwater	GEN-MW4-102921	VOAs, SVOAs	10/29/2021	11/2/2021
214958-02	Groundwater	GEN-MW1-103021	VOAs, SVOAs	10/30/2021	11/2/2021
214958-03	Groundwater	GEN-MW2-103021	VOAs, SVOAs	10/30/2021	11/2/2021
214958-04	Groundwater	GEN-DUP-103021	VOAs, SVOAs	10/30/2021	11/2/2021
214958-05	Groundwater	GEN-MW3-103121	VOAs, SVOAs	10/31/2021	11/2/2021
214958-06	Groundwater	GEN-MW6-103121	VOAs, SVOAs	10/31/2021	11/2/2021
214958-07	Groundwater	GEN-EB-110121	VOAs, SVOAs	11/1/2021	11/2/2021
214958-08	Groundwater	GEN-MW8-110121	VOAs, SVOAs	11/1/2021	11/2/2021
214958-09	Groundwater	GEN-MW7-110121	VOAs, SVOAs	11/1/2021	11/2/2021
214958-10	Water	Trip Blank T1075	VOAs	10/25/2021	11/2/2021

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Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹				
Volatile Organics	Water	Soil			
Method 8260	(µg/L)	(µg/kg) ²			
Chloromethane	5	5			
Bromomethane	5	5			
Vinyl Chloride	5	5			
Chloroethane	5	5			
Methylene Chloride	3	3			
Acetone	5	5			
Carbon Disulfide	5	5			
1,1-Dichloroethylene	5	5			
1,1-Dichloroethane	5	5			
1,2-Dichloroethylene (total)	5	5			
Chloroform	5	5			
1,2-Dichloroethane	2	2			
2-Butanone	5	5			
1,1,1-Trichloroethane	5	5			
Carbon Tetrachloride	2	2			
Bromodichloromethane	1	1			
1,2-Dichloropropane	1	1			
cis-1,3-Dichloropropene	5	5			
Trichloroethane	5	5			
Dibromochloromethane	5	5			
1,1,2-Trichloroethane	3	3			
Benzene	1	1			
trans-1,3-Dichloropropene	5	5			
Bromoform	4	4			
4-Methyl-2-pentanone	5	5			
2-Hexanone	5	5			
Tetrachloroethene	1	1			
Toluene	5	5			
1,1,2,2-Tetrachloroethane	1	1			
Chlorobenzene	5	5			
Ethylbenzene	4	4			
Styrene	5	5			
2-Chloroethyl Vinyl Ether	5	5			
1,2-Dichlorobenzene	5	5			
1,3-Dichlorobenzene	5	5			
1,4-Dichlorobenzene	5	5			
Vinyl Acetate	5	5			
Total Xylenes	5	5			
Semivolatile Organics	Water	Soil			
Method 8270	(µg/L)	(µg/kg)			
1,2,4-Trichlorobenzene	1	33			
1,2-Dichlorobenzene	10	330			
1,2-Diphenylhydrazine	10	330			
1,3-Dichlorobenzene	10	330			
1,4-Dichlorobenzene	10	330			
1,4-Dioxane	10	330			
2,4,5-Trichlorophenol	10	330			
2,4,6-Trichlorophenol	10	330			
2,4-Dichlorophenol	10	330			

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Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹			
Semivolatile Organics	Water	Soil		
Method 8270 (Cont'd.)	(µg/L)	(µg/kg)		
2,4-Dimethylphenol	10	330		
2,4-Dinitrophenol	40	1300		
2,4-Dinitrotoluene	2	67		
2,6-Dinitrotoluene	2	67		
2-Chloronaphthalene	10	330		
2-Chlorophenol	10	330		
2-Methylnaphthalene	10	330		
2-Methylphenol	10	330		
2-Nitroaniline	20	670		
2-Nitrophenol	10	330		
3,3'-Dichlorobenzidene	20	670		
3-Nitroaniline	20	670		
4,6-Dinitro-2-methylphenol	40	1300		
4-Bromophenyl-phenylether	10	330		
4-Chloro-3-methylphenol	10	330		
4-Chloroaniline	10	330		
4-Chlorophenyl-phenylether	10	330		
4-Methylphenol	10	330		
4-Nitroaniline	20	670		
4-Nitrophenol	40	1300		
Acenaphthene	10	330		
Acenaphthylene	10	330		
Acetophenone	10	330		
Aniline	10	330		
Anthracene	10	330		
Atrazine	10	330		
Benzaldehyde	10	330		
Benzidine	40	1300		
Benzo(a)anthracene	1	33		
Benzo(a)pyrene	1	33		
Benzo(b)fluoranthene	1	33		
Benzo(g,h,i)perylene	10	330		
Benzo(k)fluoranthene	1	33		
Benzoic Acid	10	330		
Benzyl Alcohol	10	330		
bis(2-chloroethoxy)methane	10	330		
bis(2-chloroethyl)ether	1	33		
bis(2-chloroisopropyl)ether	10	330		
bis(2-ethylhexyl)phthalate	10	330		
Butylbenzylphthalate	10	330		
Caprolactam	10	330		
Carbazole	10	330		
Chrysene	10	330		
Dibenzo(a,h)anthracene	1	33		
Dibenzofuran	10	330		
Diethylphthalate	10	330		
Dimethylphthalate	10	330		
Di-n-butyl phthalate	10	330		
Di-n-octyl phthalate	10	330		

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Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹		
Semivolatiles	Water	Soil	
Method 8270 (Cont'd.)	(µg/L)	(µg/kg)	
Diphenyl	10	330	
Fluoranthene	10	330	
Fluorene	10	330	
Hexachlorobenzene	1	33	
Hexachlorobutadiene	2	67	
Hexachlorocyclopentadiene	10	330	
Hexachloroethane	1	33	
Indeno(1,2,3-cd)pyrene	1	33	
Isophorone	10	330	
N,N-Dimethylaniline	1	33	
Naphthalene	10	330	
Nitrobenzene	1	33	
N-Nitrosodimethylamine	10	330	
N-Nitroso-di-n-propylamine	10	33	
	10		
N-Nitrosodiphenylamine		330	
Pentachlorophenol	40	1300	
Phenanthrene	10	330	
Phenol	10	330	
Pyrene	10	330	
Pyridine	10	330	
TAL Metals (6010/7470)	Water	Soil	
	(µg/L)	(µg/kg)	
Aluminum		40	
Antimony		2	
Arsenic		11	
Barium		40	
Beryllium		0.4	
Cadmium		1	
Calcium		1000	
Chromium		2	
Cobalt		10	
Copper		5	
Iron		30	
Lead	-	1	
Magnesium		1000	
Manganese		3	
Mercury		0.033	
Nickel		8	
Potassium	-	1000	
Selenium		1	
Silver	**	2	
Sodium	==	1000	
Thallium		2	
Vanadium		10	
Zinc		6	
Supplemental Parameters	Water	Soll	
	(µg/L)	(mg/kg)	
Total Organic Carbon (Lloyd Kahn)	NA	100	
Chloride Method 325.3	1,000	=	

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Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

Parameter	Quantitation Limit ¹			
Supplemental Parameters (Cont'd.)	Water (µg/L)	Soil (mg/kg)		
Nitrate Method 353.2	100			
Ammonia Method 350.1	100			
Iron Method 200.7	150			
Manganese Method 200.7	15			
Sulfate Method 375.4	5,000			
Sulfide Method 376.1	1,000			
Orthophosphate Method 365.2	30			
Alkalinity Method 310.1	5,000			
Methane Method 3810				
Reactive Sulfide		20		
Reactive Cyanide		25		
TCLP Benzene		1		
Total Sulfur	त्रव ा	50		
Chemical Oxygen Demand		120		

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Notes

µg/L = micrograms per liter µg/kg = micrograms per kilogram mg/kg = milligrams per kilogram

¹ Specific quantitation limits are highly matrix dependent. The quantitation limits listed are for guidance and may not always be achievable due to matrix interference.

²Quantitation limits for source materials/soil/sediment are based on wet weight. The quantitation limits calculated by the laboratory for source materials/soil/sediment (calculated on a dry weight basis) will be higher.

VOLATILE ORGANICS SAMPLE DATA



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW4-102921

Lab Sample ID:214958-01Date Sampled:10/29/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

Analyte	<u>Result</u>	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L	M	11/3/2021 13:33
Ethylbenzene	< 2.00	ug/L	M	11/3/2021 13:33
m,p-Xylene	2.44 J	ug/L		11/3/2021 13:33
o-Xylene	< 2.00	ug/L		11/3/2021 13:33
Toluene	< 2.00	ug/L	M	11/3/2021 13:33

Surrogate	Percent Recovery	<u>Limits</u>	Outliers	Date An	<u>alyzed</u>
1,2-Dichloroethane-d4	121	77.9 - 132		11/3/2021	13:33
4-Bromofluorobenzene	118	62.6 - 133		11/3/2021	13:33
Pentafluorobenzene	115	88.9 - 114	*	11/3/2021	13:33
Toluene-D8	107	75.6 - 117		11/3/2021	13:33

Method Reference(s): EPA 8260C

EPA 5030C **Data File:** z05208.D

MKP 2/16/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW1-103021

Lab Sample ID:214958-02Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 14:31
Ethylbenzene	< 2.00	ug/L		11/3/2021 14:31
m,p-Xylene	< 2.00	ug/L		11/3/2021 14:31
o-Xylene	< 2.00	ug/L		11/3/2021 14:31
Toluene	< 2.00	ug/L		11/3/2021 14:31

Surrogate	Percent Recovery	<u>Limits</u>	Outliers	Date Analyzed	
1,2-Dichloroethane-d4	117	77.9 - 132		11/3/2021	14:31
4-Bromofluorobenzene	108	62.6 - 133		11/3/2021	14:31
Pentafluorobenzene	117	88.9 - 114	*	11/3/2021	14:31
Toluene-D8	108	75.6 - 117		11/3/2021	14:31

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05211.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW2-103021

Lab Sample ID:214958-03Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

Analyte	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 14:50
Ethylbenzene	< 2.00	ug/L		11/3/2021 14:50
m,p-Xylene	1.10	ug/L	J	11/3/2021 14:50
o-Xylene	< 2.00	ug/L		11/3/2021 14:50
Toluene	< 2.00	ug/L		11/3/2021 14:50

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	<u>Outliers</u>	Date Analyzed	
1,2-Dichloroethane-d4	118	77.9 - 132		11/3/2021	14:50
4-Bromofluorobenzene	107	62.6 - 133		11/3/2021	14:50
Pentafluorobenzene	113	88.9 - 114		11/3/2021	14:50
Toluene-D8	107	75.6 - 117		11/3/2021	14:50

Method Reference(s): EPA 8260C EPA 5030C

Data File: z05212.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-DUP-103021

Lab Sample ID:214958-04Date Sampled:10/30/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 15:09
Ethylbenzene	< 2.00	ug/L		11/3/2021 15:09
m,p-Xylene	< 2.00	ug/L		11/3/2021 15:09
o-Xylene	< 2.00	ug/L		11/3/2021 15:09
Toluene	< 2.00	ug/L		11/3/2021 15:09

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	<u>Outliers</u>	Date Analyzed	
1,2-Dichloroethane-d4	123	77.9 - 132		11/3/2021	15:09
4-Bromofluorobenzene	106	62.6 - 133		11/3/2021	15:09
Pentafluorobenzene	114	88.9 - 114		11/3/2021	15:09
Toluene-D8	111	75.6 - 117		11/3/2021	15:09

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05213.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW3-103121

Lab Sample ID:214958-05Date Sampled:10/31/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 15:29
Ethylbenzene	< 2.00	ug/L		11/3/2021 15:29
m,p-Xylene	< 2.00	ug/L		11/3/2021 15:29
o-Xylene	< 2.00	ug/L		11/3/2021 15:29
Toluene	< 2.00	ug/L		11/3/2021 15:29

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	Outliers	Date Analyzed	
1,2-Dichloroethane-d4	120	77.9 - 132		11/3/2021	15:29
4-Bromofluorobenzene	118	62.6 - 133		11/3/2021	15:29
Pentafluorobenzene	110	88.9 - 114		11/3/2021	15:29
Toluene-D8	107	75.6 - 117		11/3/2021	15:29

Method Reference(s): EPA 8260C

EPA 5030C **Data File:** z05214.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW6-103121

Lab Sample ID:214958-06Date Sampled:10/31/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>	Qualifier	Date Analyzed
Benzene	178	ug/L		11/8/2021 15:52
Ethylbenzene	32.6	ug/L		11/8/2021 15:52
m,p-Xylene	43.5	ug/L		11/8/2021 15:52
o-Xylene	43.3	ug/L		11/8/2021 15:52
Toluene	33.8	ug/L		11/8/2021 15:52

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	Outliers	Date Analyzed	
1,2-Dichloroethane-d4	99.0	77.9 - 132		11/8/2021	15:52
4-Bromofluorobenzene	86.5	62.6 - 133		11/8/2021	15:52
Pentafluorobenzene	99.3	88.9 - 114		11/8/2021	15:52
Toluene-D8	88.7	75.6 - 117		11/8/2021	15:52

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05290.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-EB-110121

Lab Sample ID:214958-07Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	0.806 J	ug/L	J	11/3/2021 16:07
Ethylbenzene	< 2.00	ug/L		11/3/2021 16:07
m,p-Xylene	< 2.00	ug/L		11/3/2021 16:07
o-Xylene	< 2.00	ug/L		11/3/2021 16:07
Toluene	< 2.00	ug/L		11/3/2021 16:07

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	Outliers	<u>S Date Analyzed</u>	
1,2-Dichloroethane-d4	117	77.9 - 132		11/3/2021	16:07
4-Bromofluorobenzene	117	62.6 - 133		11/3/2021	16:07
Pentafluorobenzene	116	88.9 - 114	*	11/3/2021	16:07
Toluene-D8	108	75.6 - 117		11/3/2021	16:07

Method Reference(s): EPA 8260C EPA 5030C

Data File: z05216.D

MKP 2/16/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW8-110121

Lab Sample ID:214958-08Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>	Qualifier	Date Analyzed
Benzene	3.37 J	ug/L		11/3/2021 16:27
Ethylbenzene	2.32 J	ug/L		11/3/2021 16:27
m,p-Xylene	3.52 J	ug/L		11/3/2021 16:27
o-Xylene	3.06 J	ug/L		11/3/2021 16:27
Toluene	1.44 J	ug/L	J	11/3/2021 16:27

<u>Surrogate</u>	Percent Recovery	<u>very Limits Outliers</u>		<u>Date An</u>	<u>alyzed</u>
1,2-Dichloroethane-d4	118	77.9 - 132		11/3/2021	16:27
4-Bromofluorobenzene	112	62.6 - 133		11/3/2021	16:27
Pentafluorobenzene	115	88.9 - 114	*	11/3/2021	16:27
Toluene-D8	97.0	75.6 - 117		11/3/2021	16:27

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05217.D

MKP 2/16/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN-MW7-110121

Lab Sample ID:214958-09Date Sampled:11/1/2021Matrix:GroundwaterDate Received:11/2/2021

Volatile Organics

Analyte	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 16:46
Ethylbenzene	< 2.00	ug/L		11/3/2021 16:46
m,p-Xylene	< 2.00	ug/L		11/3/2021 16:46
o-Xylene	< 2.00	ug/L		11/3/2021 16:46
Toluene	< 2.00	ug/L		11/3/2021 16:46

Surrogate	Percent Recovery	<u>Limits</u>	<u>Outliers</u>	Date Analyzed	
1,2-Dichloroethane-d4	115	77.9 - 132		11/3/2021	16:46
4-Bromofluorobenzene	101	62.6 - 133		11/3/2021	16:46
Pentafluorobenzene	114	88.9 - 114		11/3/2021	16:46
Toluene-D8	97.7	75.6 - 117		11/3/2021	16:46

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z05218.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: Trip Blank T1075

 Lab Sample ID:
 214958-10
 Date Sampled:
 10/25/2021

 Matrix:
 Water
 Date Received:
 11/2/2021

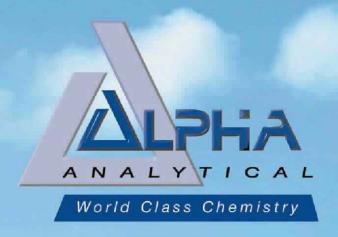
Volatile Organics

Analyte	<u>Result</u>	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		11/3/2021 17:05
Ethylbenzene	< 2.00	ug/L		11/3/2021 17:05
m,p-Xylene	< 2.00	ug/L		11/3/2021 17:05
o-Xylene	< 2.00	ug/L		11/3/2021 17:05
Toluene	< 2.00	ug/L		11/3/2021 17:05

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	<u>Outliers</u>	Date Analyzed		
1,2-Dichloroethane-d4	127	77.9 - 132		11/3/2021	17:05	
4-Bromofluorobenzene	111	62.6 - 133		11/3/2021	17:05	
Pentafluorobenzene	117	88.9 - 114	*	11/3/2021	17:05	
Toluene-D8	121	75.6 - 117	*	11/3/2021	17:05	

Method Reference(s): EPA 8260C

EPA 5030C **Data File:** z05219.D



www.alphalab.com



Alpha Analytical

Laboratory Code: 11148

SDG Number: L2160069

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 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2160069-01	GEN-MW4-102921 214958-01	WATER	Not Specified	10/29/21 16:15	11/02/21
L2160069-02	GEN-MW1-103021 214958-02	WATER	Not Specified	10/30/21 14:00	11/02/21
L2160069-03	GEN-MW2-103021 214958-03	WATER	Not Specified	10/30/21 15:45	11/02/21
L2160069-04	GEN-DUP-103021 214958-04	WATER	Not Specified	10/30/21 00:00	11/02/21
L2160069-05	GEN-MW3-103121 214958-05	WATER	Not Specified	10/31/21 12:30	11/02/21
L2160069-06	GEN-MW6-103121 214958-06	WATER	Not Specified	10/31/21 14:30	11/02/21
L2160069-07	GEN-EB-110121 214958-07	WATER	Not Specified	11/01/21 12:00	11/02/21
L2160069-08	GEN-MW8-110121 214958-08	WATER	Not Specified	11/01/21 15:00	11/02/21
L2160069-09	GEN-MW7-110121 214958-09	WATER	Not Specified	11/01/21 16:30	11/02/21



 Project Name:
 214958
 Lab Number:
 L2160069

 Project Number:
 214958
 Report Date:
 11/09/21

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature: (attlin Wallieru

Report Date: 11/09/21

Title: Technical Director/Representative

INVOICE TO:

Lallebole9

11148



CHAIN OF CUSTODY

REPORT TO:

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Semivolatiles Data by Method 8270D-SIM

No data validation qualifiers were added

MKP 2/16/2022

Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-01

Client ID : GEN-MW4-102921 214958-01

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-01
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 10/29/21 16:15

Date Received : 11/02/21

Date Analyzed : 11/04/21 20:06

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : ALS
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
83-32-9	Acenaphthene	ND	0.10	0.04	U
206-44-0	Fluoranthene	ND	0.10	0.04	U
91-20-3	Naphthalene	ND	0.10	0.04	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U
218-01-9	Chrysene	ND	0.10	0.04	U
208-96-8	Acenaphthylene	ND	0.10	0.04	U
120-12-7	Anthracene	ND	0.10	0.04	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U
86-73-7	Fluorene	ND	0.10	0.04	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U
129-00-0	Pyrene	ND	0.10	0.04	U



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-02

Client ID : GEN-MW1-103021 214958-02

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-02
Sample Amount : 1000 ml
Extraction Method : EPA 3510C

Extract Volume : 1000 uL GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 10/30/21 14:00

Date Received : 11/02/21

Date Analyzed : 11/05/21 17:19

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L				
CAS NO.	Parameter	Results	RL	MDL	Qualifier		
83-32-9	Acenaphthene	ND	0.10	0.04	U		
206-44-0	Fluoranthene	ND	0.10	0.04	U		
91-20-3	Naphthalene	ND	0.10	0.04	U		
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U		
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U		
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U		
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U		
218-01-9	Chrysene	ND	0.10	0.04	U		
208-96-8	Acenaphthylene	ND	0.10	0.04	U		
120-12-7	Anthracene	ND	0.10	0.04	U		
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U		
86-73-7	Fluorene	ND	0.10	0.04	U		
85-01-8	Phenanthrene	ND	0.10	0.02	U		
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U		
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U		
129-00-0	Pyrene	ND	0.10	0.04	U		



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-03

Client ID : GEN-MW2-103021 214958-03

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-03
Sample Amount : 1000 ml
Extraction Method : EPA 3510C

Extract Volume : 1000 uL GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 10/30/21 15:45

Date Received : 11/02/21

Date Analyzed : 11/05/21 17:35

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

%Solids : N/A Injection Volume : 1 uL

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	0.02	0.10	0.02	J	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-04

Client ID : GEN-DUP-103021 214958-04

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-04
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 10/30/21 00:00

Date Received : 11/02/21

Date Analyzed : 11/05/21 17:52

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L				
CAS NO.	Parameter	Results	RL	MDL	Qualifier		
83-32-9	Acenaphthene	ND	0.10	0.04	U		
206-44-0	Fluoranthene	ND	0.10	0.04	U		
91-20-3	Naphthalene	ND	0.10	0.04	U		
56-55-3	Benzo(a)anthracene	0.02	0.10	0.02	J		
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U		
205-99-2	Benzo(b)fluoranthene	0.03	0.10	0.02	J		
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U		
218-01-9	Chrysene	ND	0.10	0.04	U		
208-96-8	Acenaphthylene	ND	0.10	0.04	U		
120-12-7	Anthracene	ND	0.10	0.04	U		
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U		
86-73-7	Fluorene	ND	0.10	0.04	U		
85-01-8	Phenanthrene	0.02	0.10	0.02	J		
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U		
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U		
129-00-0	Pyrene	ND	0.10	0.04	U		



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-05

Client ID : GEN-MW3-103121 214958-05

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-05
Sample Amount : 1000 ml

Extraction Method : EPA 3510C Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 10/31/21 12:30

Date Received : 11/02/21

Date Analyzed : 11/05/21 18:08

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	



Client : Paradigm Environmental Services

Project Name : 214958

Lab ID : L2160069-06D

Client ID : GEN-MW6-103121 214958-06

Sample Location :

GPC Cleanup

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 069-06D1
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

: 1000 uL : N

%Solids : N/A Injection Volume : 1 uL

Lab Number

Date Analyzed

Date Extracted

Dilution Factor

Instrument ID

GC Column

Analyst

Project Number : 214958

Date Received : 11/02/21

Date Collected : 10/31/21 14:30

: L2160069

: 11/03/21

: SV115

: RXI-5SilM

: 10

: 11/09/21 11:40

		ug/L				
Parameter	Results	RL	MDL	Qualifier		
Acenaphthene	1.7	1.0	0.35			
Fluoranthene	ND	1.0	0.38	U		
Naphthalene	170	1.0	0.43			
Benzo(a)anthracene	ND	1.0	0.18	U		
Benzo(a)pyrene	ND	1.0	0.39	U		
Benzo(b)fluoranthene	ND	1.0	0.16	U		
Benzo(k)fluoranthene	ND	1.0	0.42	U		
Chrysene	ND	1.0	0.38	U		
Acenaphthylene	18	1.0	0.35			
Anthracene	0.46	1.0	0.35	J		
Benzo(ghi)perylene	ND	1.0	0.42	U		
Fluorene	4.0	1.0	0.37			
Phenanthrene	1.6	1.0	0.15			
Dibenzo(a,h)anthracene	ND	1.0	0.39	U		
Indeno(1,2,3-cd)pyrene	ND	1.0	0.40	U		
Pyrene	ND	1.0	0.40	U		
	Acenaphthene Fluoranthene Naphthalene Benzo(a)anthracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(k)fluoranthene Chrysene Acenaphthylene Anthracene Benzo(ghi)perylene Fluorene Phenanthrene Dibenzo(a,h)anthracene Indeno(1,2,3-cd)pyrene	Acenaphthene 1.7 Fluoranthene ND Naphthalene 170 Benzo(a)anthracene ND Benzo(a)pyrene ND Benzo(b)fluoranthene ND Benzo(k)fluoranthene ND Chrysene ND Acenaphthylene 18 Anthracene 0.46 Benzo(ghi)perylene ND Fluorene 4.0 Phenanthrene 1.6 Dibenzo(a,h)anthracene ND Indeno(1,2,3-cd)pyrene ND	Parameter Results RL Acenaphthene 1.7 1.0 Fluoranthene ND 1.0 Naphthalene 170 1.0 Benzo(a)anthracene ND 1.0 Benzo(b)fluoranthene ND 1.0 Benzo(k)fluoranthene ND 1.0 Chrysene ND 1.0 Acenaphthylene 18 1.0 Anthracene 0.46 1.0 Benzo(ghi)perylene ND 1.0 Fluorene 4.0 1.0 Phenanthrene 1.6 1.0 Dibenzo(a,h)anthracene ND 1.0 Indeno(1,2,3-cd)pyrene ND 1.0	Parameter Results RL MDL Acenaphthene 1.7 1.0 0.35 Fluoranthene ND 1.0 0.38 Naphthalene 170 1.0 0.43 Benzo(a)anthracene ND 1.0 0.18 Benzo(a)pyrene ND 1.0 0.39 Benzo(b)fluoranthene ND 1.0 0.16 Benzo(k)fluoranthene ND 1.0 0.42 Chrysene ND 1.0 0.38 Acenaphthylene 18 1.0 0.35 Anthracene 0.46 1.0 0.35 Benzo(ghi)perylene ND 1.0 0.42 Fluorene 4.0 1.0 0.37 Phenanthrene 1.6 1.0 0.15 Dibenzo(a,h)anthracene ND 1.0 0.39 Indeno(1,2,3-cd)pyrene ND 1.0 0.40	Parameter Results RL MDL Qualifier Acenaphthene 1.7 1.0 0.35 Fluoranthene ND 1.0 0.38 U Naphthalene 170 1.0 0.43 U Benzo(a)anthracene ND 1.0 0.18 U Benzo(a)pyrene ND 1.0 0.39 U Benzo(b)fluoranthene ND 1.0 0.16 U Benzo(k)fluoranthene ND 1.0 0.42 U Chrysene ND 1.0 0.38 U Acenaphthylene 18 1.0 0.35 J Anthracene 0.46 1.0 0.35 J Benzo(ghi)perylene ND 1.0 0.37 U Fluorene 4.0 1.0 0.37 U Phenanthrene 1.6 1.0 0.15 Dibenzo(a,h)anthracene ND 1.0 0.40 U	



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-07

Client ID : GEN-EB-110121 214958-07

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-07
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 11/01/21 12:00

Date Received : 11/02/21

Date Analyzed : 11/05/21 18:41

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	



Client : Paradigm Environmental Services

Project Name : 214958 Lab ID : L2160069-08

Client ID : GEN-MW8-110121 214958-08

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 60069-08
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2160069 Project Number : 214958

Date Collected : 11/01/21 15:00

Date Received : 11/02/21

Date Analyzed : 11/05/21 18:58

Date Extracted : 11/03/21

Dilution Factor : 1
Analyst : JJW
Instrument ID : SV120
GC Column : RXI-5SiIM
%Solids : N/A

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	0.22	0.10	0.04		
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	0.04	0.10	0.02	J	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	



Client : Paradigm Environmental Services

Project Name : 214958 : L2160069-09 Lab ID

Client ID : GEN-MW7-110121 214958-09

Sample Location :

Sample Matrix : WATER Analytical Method : 1,8270D-SIM Lab File ID : 60069-09 Sample Amount : 1000 ml Extraction Method : EPA 3510C **Extract Volume** : 1000 uL

GPC Cleanup : N Lab Number : L2160069 Project Number : 214958

Date Collected : 11/01/21 16:30 Date Received : 11/02/21

Date Analyzed : 11/05/21 19:14

Date Extracted : 11/03/21

Dilution Factor Analyst : JJW Instrument ID : SV120 GC Column : RXI-5SilM %Solids : N/A

			ug/L				
CAS NO.	Parameter	Results	RL	MDL	Qualifier	_	
83-32-9	Acenaphthene	ND	0.10	0.04	U		
206-44-0	Fluoranthene	ND	0.10	0.04	U		
91-20-3	Naphthalene	ND	0.10	0.04	U		
56-55-3	Benzo(a)anthracene	0.02	0.10	0.02	J		
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U		
205-99-2	Benzo(b)fluoranthene	0.02	0.10	0.02	J		
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U		
218-01-9	Chrysene	ND	0.10	0.04	U		
208-96-8	Acenaphthylene	ND	0.10	0.04	U		
120-12-7	Anthracene	ND	0.10	0.04	U		
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U		
86-73-7	Fluorene	ND	0.10	0.04	U		
35-01-8	Phenanthrene	0.02	0.10	0.02	J		
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U		
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U		
129-00-0	Pyrene	ND	0.10	0.04	U		



Appendix B

Laboratory QC Documentation

2 VOLATILE SURROGATE RECOVERY

Lab Name: <u>Paradigm Environmental Services</u>

Lab Project #: 214958
Client Name: Neu-Velle

Client Project Name: RGE Geneseo Fmr. MGP Site

Client Project #: N/A

SDG No.: 4958-01 Matrix: Groundwater
QC Batch: QC211103VOAW

Instrument ID: <u>Instrument1</u>

GC Column 1: $\underline{DB-624}$ ID (mm): $\underline{0.20}$ Detector: \underline{MSD}

LAB SAMPLE NO.	CLIENT SAMPLE ID	PFB %REC	12DCEd4 %REC	TD8 %REC	4BFB %REC	Total Out
1 Blk 1	N/A	116 *	119	118 *	111	2
2 LCS 1	N/A	111	116	126 *	109	1
3 214958-01	GEN-MW4-102921	115 *	121	107	118	1
4 214958-01MS	GEN-MW4-102921	109	114	123 *	92.6	1
5 214958-01MSD	GEN-MW4-102921	110	113	116	89.9	0
6 214958-02	GEN-MW1-103021	117 *	117	108	108	1
7 214958-03	GEN-MW2-103021	113	118	107	107	0
8 214958-04	GEN-DUP-103021	114	123	111	106	0
9 214958-05	GEN-MW3-103121	110	120	107	118	0
10 214958-07	GEN-EB-110121	116 *	117	108	117	1
11 214958-08	GEN-MW8-110121	115 *	118	97.0	112	1
12 214958-09	GEN-MW7-110121	114	115	97.7	101	0
13 214958-10	Trip Blank T1075	117 *	127	121 *	111	2
14						
15						
16						
17						
18						
19						
20						
21						
22						
23		-				
24		-				
25						

	QC LIMITS %
PFB = Pentafluorobenzene	(88.9 - 114)
12DCEd4 = 1,2-Dichloroethane-d4	(77.9 - 132)
TD8 = Toluene-d8	(75.6 - 117)
4BFB = 4-Bromofluorobenzene	(62.6 - 133)

^{*} Values outside of current required QC limits

D Surrogate diluted out



QC Report for Matrix Spike and Matrix Spike Duplicate

SDG #: 4958-01

Neu-Velle Lab Project ID: 214958

Project Reference: RGE Geneseo Fmr. MGP Site

Lab Sample ID: 214958-01

Sample Identifier: GEN-MW4-102921

Matrix: Groundwater

Date Sampled: 10/29/2021 **Date Received:** 11/2/2021

Date Analyzed: 11/3/2021

Volatile Organics

Client:

	Sample	Result	<u>MS</u>	<u>MS</u>	MS %	MSD	MSD	MSD %	<u>% Rec.</u>	<u>MS</u>	<u>MSD</u>	<u>Relative</u>	<u>RPD</u>	<u>RPD</u>
<u>Analyte</u>	Result	<u>Units</u>	Added	Result	Recovery	Added	Result	Recovery	<u>Limits</u>	<u>Outlier</u>	<u>Outlier</u>	% Diff.	Limit	<u>Outlier</u>
Benzene	< 1.00	ug/L	50.0	60.0	120	50.0	62.3	125	85.6 - 106	*	*	3.75	10.2	
Ethylbenzene	< 2.00	ug/L	50.0	53.3	107	50.0	54.8	110	80.5 - 106	*	*	2.85	15.3	
Toluene	< 2.00	ug/L	50.0	67.8	136	50.0	66.6	133	72.9 - 107	*	*	1.77	18.1	

Method Reference(s): EPA 8260C

EPA 5030C

Data File(s): z05209.D

z05210.D z05208.D

QC Batch ID: QC211103V0AW

Any estimated values are displayed, and derived values calculated, based on numeric result only. See primary analytical report for data flags.

Appendix C

Validator Qualifications

KENNETH R. APPLIN Geochemist/Data Validator

Ph.D., Geochemistry and Mineralogy, The Pennsylvania State University

M.S., Geochemistry and Mineralogy, The Pennsylvania State University

B.A., Geological Sciences, SUNY at Geneseo, NY

Dr. Applin has over 35 years of experience working with the geochemistry of natural waters. His prior experience includes working as an Assistant Professor of Geology at the University of Missouri-Columbia and as Chief Hydrogeologist and Geochemist with a leading engineering firm in Rochester, NY. In 1993, he established KR Applin and Associates, a small consulting business that focuses on the geochemistry of natural waters, especially as applied to problems involving the contamination of groundwater and surface water.

Dr. Applin is also an experienced analytical data validator and has provided data validation services since 1994 to a variety of clients performing brownfield cleanup projects, hazardous waste remediation, groundwater monitoring at solid waste facilities, and other projects requiring third-party data validation. Dr. Applin has several years of hands-on experience with the laboratory analysis of natural waters and has successfully completed the USEPA Region II certification courses for performing inorganic and organic analytical data validation.

MICHAEL K. PERRY Chemist/Data Validator

B.S. Chemistry, Georgia State University, Atlanta, GA

A.A.S., Chemical Technology, Alfred State College, Alfred, NY

Mr. Perry has over 30 years of experience in the analytical laboratory business. During his early career, he spent several years as a laboratory analyst performing the analysis of soil, water, and air samples for inorganic and organic chemical parameters. During his last 20 years in the environmental laboratory business, he managed and directed two major analytical laboratories in Rochester, NY. His management responsibilities included oversight of the daily operations of the lab, staff training and supervision, the selection, purchase, and maintenance of analytical instruments, the introduction of new laboratory methods, analytical quality assurance and quality control, data acquisition and management, and other business-related activities.

Mr. Perry has an extensive working knowledge of the methods and procedures used for sampling and analyzing both inorganic and organic analytes in soil, water, and air. He is an accomplished laboratory chemist and is familiar with the analytical methods and procedures established under the USEPA Contract Laboratory Protocols (CLP), the NYSDEC Analytical Services Protocols (ASP), and the NYSDOH Environmental Laboratory Approval Program (ELAP).



Analytical Report For

Neu-Velle

For Lab Project ID

222457

Referencing

RGE Geneseo Fmr. MGP Site

Prepared

Friday, June 10, 2022

Any noncompliant QC parameters or other notes impacting data interpretation are flagged or documented on the final report or are noted below:

Portions of the enclosed report reflects analysis that has been subcontracted and are presented in their original form.

Emily Farmen

Certifies that this report has been approved by the Technical Director or Designee

179 Lake Avenue • Rochester, NY 14608 • (585) 647-2530 • Fax (585) 647-3311 • ELAP ID# 10958



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW3 - 052322

Lab Sample ID: 222457-01 **Date Sampled:** 5/23/2022 14:00

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Result</u>	<u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
< 1.00	ug/L			6/2/20	22 13:30
< 2.00	ug/L			6/2/20	22 13:30
< 2.00	ug/L			6/2/20	22 13:30
< 2.00	ug/L			6/2/20	22 13:30
< 2.00	ug/L			6/2/20	22 13:30
Perce	ent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
	83.7	81.1 - 136		6/2/2022	13:30
	87.0	75.8 - 132		6/2/2022	13:30
	121	82 - 132		6/2/2022	13:30
	115	64.6 - 137		6/2/2022	13:30
	< 1.00 < 2.00 < 2.00 < 2.00 < 2.00 < 2.00	< 1.00 ug/L < 2.00 ug/L Percent Recovery 83.7 87.0 121	< 1.00 ug/L < 2.00 ug/L Percent Recovery 83.7 81.1 - 136 87.0 75.8 - 132 121 82 - 132	< 1.00	< 1.00

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09594.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - EB - 052322

Lab Sample ID: 222457-02 **Date Sampled:** 5/23/2022 12:45

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date An	alyzed
Benzene	< 1.00	ug/L			6/2/202	22 13:49
Ethylbenzene	< 2.00	ug/L			6/2/202	22 13:49
m,p-Xylene	< 2.00	ug/L			6/2/202	22 13:49
o-Xylene	< 2.00	ug/L			6/2/202	22 13:49
Toluene	< 2.00	ug/L			6/2/202	22 13:49
<u>Surrogate</u>	Perce	nt Recovery	<u>Limits</u>	Outliers	Date Ana	alyzed
1,2-Dichloroethane-d4		74.1	81.1 - 136	*	6/2/2022	13:49
4-Bromofluorobenzene		76.6	75.8 - 132		6/2/2022	13:49
Pentafluorobenzene		126	82 - 132		6/2/2022	13:49
Toluene-D8		106	64.6 - 137		6/2/2022	13:49

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09595.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - DUP - 052322

Lab Sample ID:222457-03Date Sampled: 5/23/2022Matrix:GroundwaterDate Received 5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
Benzene	< 1.00	ug/L			6/2/20	22 14:09
Ethylbenzene	< 2.00	ug/L			6/2/20	22 14:09
m,p-Xylene	< 2.00	ug/L			6/2/20	22 14:09
o-Xylene	< 2.00	ug/L			6/2/20	22 14:09
Toluene	< 2.00	ug/L			6/2/20	22 14:09
Surrogate	Perc	ent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		72.5	81.1 - 136	*	6/2/2022	14:09
4-Bromofluorobenzene		74.3	75.8 - 132	*	6/2/2022	14:09
Pentafluorobenzene		129	82 - 132		6/2/2022	14:09
Toluene-D8		109	64.6 - 137		6/2/2022	14:09

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09596.D



Client: Neu-Velle

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW2 - 052322

Lab Sample ID: 222457-04 **Date Sampled:** 5/23/2022 15:30

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>		Qualifier	Date An	alyzed
Benzene	< 1.00	ug/L			6/2/202	22 14:28
Ethylbenzene	< 2.00	ug/L		M	6/2/202	22 14:28
m,p-Xylene	< 2.00	ug/L			6/2/202	22 14:28
o-Xylene	< 2.00	ug/L			6/2/202	22 14:28
Toluene	< 2.00	ug/L			6/2/202	22 14:28
Surrogate	<u>Perce</u>	nt Recovery	<u>Limits</u>	Outliers	Date Ana	alyzed
1,2-Dichloroethane-d4		70.5	81.1 - 136	*	6/2/2022	14:28
4-Bromofluorobenzene		68.1	75.8 - 132	*	6/2/2022	14:28
Pentafluorobenzene		126	82 - 132		6/2/2022	14:28
Toluene-D8		100	64.6 - 137		6/2/2022	14:28

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09597.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW1 - 052422

Lab Sample ID: 222457-05 **Date Sampled:** 5/24/2022 14:45

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

Analyte	<u>Result</u>	<u>Units</u>		Qualifier	Date Ar	<u>nalyzed</u>
Benzene	< 1.00	ug/L			6/2/20	22 15:45
Ethylbenzene	< 2.00	ug/L			6/2/20	22 15:45
m,p-Xylene	< 2.00	ug/L			6/2/20	22 15:45
o-Xylene	< 2.00	ug/L			6/2/20	22 15:45
Toluene	< 2.00	ug/L			6/2/20	22 15:45
Surrogate	Perc	ent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		72.8	81.1 - 136	*	6/2/2022	15:45
4-Bromofluorobenzene		72.2	75.8 - 132	*	6/2/2022	15:45
Pentafluorobenzene		125	82 - 132		6/2/2022	15:45
Toluene-D8		106	64.6 - 137		6/2/2022	15:45

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09601.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW4 - 052422

Lab Sample ID: 222457-06 **Date Sampled:** 5/24/2022 16:30

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date An	<u>alyzed</u>
Benzene	0.503	ug/L		J	6/2/202	2 16:05
Ethylbenzene	< 2.00	ug/L			6/2/202	2 16:05
m,p-Xylene	2.98	ug/L			6/2/202	2 16:05
o-Xylene	< 2.00	ug/L			6/2/202	2 16:05
Toluene	< 2.00	ug/L			6/2/202	2 16:05
<u>Surrogate</u>	Percent	Recovery	<u>Limits</u>	Outliers	Date Ana	lyzed
1,2-Dichloroethane-d4	6	8.6	81.1 - 136	*	6/2/2022	16:05
4-Bromofluorobenzene	68	8.9	75.8 - 132	*	6/2/2022	16:05
Pentafluorobenzene	1	22	82 - 132		6/2/2022	16:05
Toluene-D8	98	8.6	64.6 - 137		6/2/2022	16:05

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09602.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW6 - 052522

Lab Sample ID: 222457-07 **Date Sampled:** 5/25/2022 14:45

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

Analyte	Res	<u>sult</u> <u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
Benzene	164	ug/L			6/2/20	22 15:26
Ethylbenzene	46.7	ug/L			6/2/20	22 15:26
m,p-Xylene	55.7	ug/L			6/2/20	22 15:26
o-Xylene	48.5	ug/L			6/2/20	22 15:26
Toluene	49.8	ug/L			6/2/20	22 15:26
<u>Surrogate</u>		Percent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		79.8	81.1 - 136	*	6/2/2022	15:26
4-Bromofluorobenzene		88.8	75.8 - 132		6/2/2022	15:26
Pentafluorobenzene		122	82 - 132		6/2/2022	15:26
Toluene-D8		108	64.6 - 137		6/2/2022	15:26

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09600.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW8 - 052522

Lab Sample ID: 222457-08 **Date Sampled:** 5/25/2022 16:30

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	Res	sult <u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
Benzene	1.88	ug/L			6/2/202	22 16:24
Ethylbenzene	3.08	ug/L			6/2/202	22 16:24
m,p-Xylene	5.86	ug/L			6/2/202	22 16:24
o-Xylene	3.12	ug/L			6/2/202	22 16:24
Toluene	2.00	ug/L		J	6/2/202	22 16:24
<u>Surrogate</u>		Percent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		75.4	81.1 - 136	*	6/2/2022	16:24
4-Bromofluorobenzene		79.3	75.8 - 132		6/2/2022	16:24
Pentafluorobenzene		121	82 - 132		6/2/2022	16:24
Toluene-D8		106	64.6 - 137		6/2/2022	16:24

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09603.D



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW7 - 052622

Lab Sample ID: 222457-09 **Date Sampled:** 5/26/2022 13:00

Matrix: Groundwater Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
Benzene	< 1.00	ug/L			6/2/20	22 16:43
Ethylbenzene	< 2.00	ug/L			6/2/20	22 16:43
m,p-Xylene	< 2.00	ug/L			6/2/20	22 16:43
o-Xylene	< 2.00	ug/L			6/2/20	22 16:43
Toluene	< 2.00	ug/L			6/2/20	22 16:43
Surrogate	Perc	ent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		81.6	81.1 - 136		6/2/2022	16:43
4-Bromofluorobenzene		78.8	75.8 - 132		6/2/2022	16:43
Pentafluorobenzene		113	82 - 132		6/2/2022	16:43
Toluene-D8		103	64.6 - 137		6/2/2022	16:43

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09604.D



Lab Project ID: 222457

Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: Trip Blank

 Lab Sample ID:
 222457-10
 Date Sampled: 5/26/2022

 Matrix:
 Water
 Date Received 5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>		Qualifier	<u>Date An</u>	<u>alyzed</u>
Benzene	< 1.00	ug/L			6/2/202	22 12:07
Ethylbenzene	< 2.00	ug/L			6/2/202	22 12:07
m,p-Xylene	< 2.00	ug/L			6/2/202	22 12:07
o-Xylene	< 2.00	ug/L			6/2/202	22 12:07
Toluene	< 2.00	ug/L			6/2/202	22 12:07
<u>Surrogate</u>	Percen	t Recovery	<u>Limits</u>	Outliers	Date Ana	llyzed
1,2-Dichloroethane-d4	7	75.8	81.1 - 136	*	6/2/2022	12:07
4-Bromofluorobenzene	7	78.7	75.8 - 132		6/2/2022	12:07
Pentafluorobenzene	2	121	82 - 132		6/2/2022	12:07
Toluene-D8	2	107	64.6 - 137		6/2/2022	12:07

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09590.D



Analytical Report Appendix

The reported results relate only to the samples as they have been received by the laboratory.

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All soil/sludge samples have been reported on a dry weight basis, unless qualified "reported as received". Other solids are reported as received.

Low level Volatiles blank reports for soil/solid matrix are based on a nominal 5 gram weight. Sample results and reporting limits are based on actual weight, which may be more or less than 5 grams.

The Chain of Custody provides additional information, including compliance with sample condition requirements upon receipt. Sample condition requirements are defined under the 2003 NELAC Standard, sections 5.5.8.3.1 and 5.5.8.3.2.

NYSDOH ELAP does not certify for all parameters. Paradigm Environmental Services or the indicated subcontracted laboratory does hold certification for all analytes where certification is offered by ELAP unless otherwise specified. Aliquots separated for certain tests, such as TCLP, are indicated on the Chain of Custody and final reports with an "A" suffix.

Data qualifiers are used, when necessary, to provide additional information about the data. This information may be communicated as a flag or as text at the bottom of the report. Please refer to the following list of analyte-specific, frequently used data flags and their meaning:

- "<" = Analyzed for but not detected at or above the quantitation limit.
- "E" = Result has been estimated, calibration limit exceeded.
- "H" = Denotes a parameter analyzed outside of holding time.
- "Z" = See case narrative.
- "D" = Sample, Laboratory Control Sample, or Matrix Spike Duplicate results above Relative Percent Difference limit.
- "M" = Matrix spike recoveries outside QC limits. Matrix bias indicated.
- "B" = Method blank contained trace levels of analyte. Refer to included method blank report.
- "J" = Result estimated between the quantitation limit and half the quantitation limit.
- "L" = Laboratory Control Sample recovery outside accepted QC limits.
- "P" = Concentration differs by more than 40% between the primary and secondary analytical columns.
- "NC" = Not calculable. Applicable to RPD if sample or duplicate result is non-detect or estimated (see primary report for data flags). Applicable to MS if sample is greater or equal to ten times the spike added. Applicable to sample surrogates or MS if sample dilution is 10x or higher.
- "*" = Indicates any recoveries outside associated acceptance windows. Surrogate outliers in samples are presumed matrix effects. LCS demonstrates method compliance unless otherwise noted.
- "(1)" = Indicates data from primary column used for QC calculation.
- "A" = denotes a parameter for which ELAP does not offer approval as part of their laboratory certification program.
- "F" = denotes a parameter for which Paradigm does not carry certification, the results for which should therefore only be used where ELAP certification is not required, such as personal exposure assessment.

This report is part of a multipage document and should only be evaluated in its entirety. The Chain of Custody provides additional sample information, including compliance with the sample condition requirements upon receipt.

GENERAL TERMS AND CONDITIONS LABORATORY SERVICES

These Terms and Conditions embody the whole agreement of the parties in the absence of a signed and executed contract between the Laboratory (LAB) and Client. They shall supersede all previous communications, representations, or agreements, either verbal or written, between the parties. The LAB specifically rejects all additional, inconsistent, or conflicting terms, whether printed or otherwise set forth in any purchase order or other communication from the Client to the LAB. The invalidity or unenforceability in whole or in part of any provision, tern or condition hereof shall not affect in any way the validity or enforceability of the remainder of the Terms and Conditions. No waiver by LAB of any provision, term, or condition hereof or of any breach by or obligation of the Client hereunder shall constitute a waiver of such provision, term, or condition on any other occasion or a waiver of any other breach by or obligation of the Client. This agreement shall be administered and interpreted under the laws of the state which services are procured.

Warranty.

Recognizing that the nature of many samples is unknown and that some may contain potentially hazardous components, LAB warrants only that it will perform testing services, obtain findings, and prepare reports in accordance with generally accepted analytical laboratory principles and practices at the time of performance of services. LAB makes no other warranty, express or implied.

Scope and Compensation. LAB agrees to perform the services described in the chain of custody to which these terms and conditions are attached. Unless the parties agree in writing to the contrary, the duties of LAB shall not be construed to exceed the services specifically described. LAB wi use LAB default method for all tests unless specified otherwise on the Work Order.

Payment terms are net 30 days from the date of invoice. All overdue payments are subject to an interest charge of one and one-half percent (1-1/2%) per month or a portion thereof. Client shall also be responsible for costs of collection, including payment of reasonable attorney fees if such expense is incurred. The prices, unless stated, do not include any sale, use or other taxes. Such taxes will be added to invoice prices when required.

Prices.

Compensation for services performed will be based on the current Lab Analytical Fee Schedule or on quotations agreed to in writing by the parties. Turnaround time based charges are determined from the time of resolution of all work order questions. Testimony, court appearances or data compilation for legal action will be charged separately. Evaluation and reporting of initial screening runs may incur additional fees.

Limitations of Liability.

In the event of any error, omission, or other professional negligence, the sole and exclusive responsibility of LAB shall be to reperform the deficient work at its own expense and LAB shall have no other liability whatsoever. All claims shall be deemed waived unless made in writing and received by LAB within ninety (90) days following completion of services.

LAB shall have no liability, obligation, or responsibility of any kind for losses, costs, expenses, or other damages (including but not limited to any special, direct, incidental or consequential damages) with respect to LAB's services or results.

All results provided by LAB are strictly for the use of its clients and LAB is in no way responsible for the use of such results by clients or third parties. All reports should be considered in their entirety, and LAB is not responsible for the separation, detachment, or other use of any portion of these reports. Client may not assign the lab report without the written consent of the LAB. Client covenants and agrees, at its/his/her sole expense, to indemnify, protect, defend, and save harmless the LAB from and against

any and all damages, losses, liabilities, obligations, penalties, claims, litigation, demands, defenses, judgments, suits, actions, proceedings, costs, disbursements and/or expenses (including, without limitation attorneys' and experts' fees and disbursements) of any kind whatsoever which may at any time be imposed upon, incurred by or asserted or awarded against client relating to, resulting from or arising out of (a) the breach of this agreement by this client, (b) the negligence of the client in handling, delivering or disclosing any hazardous substance, (c) the violation of the Client of any applicable law, (d) non-compliance by the Client with any

environmental permit or (e) a material misrepresentation in disclosing the materials to be tested.

Hazard Disclosure.

Client represents and warrants that any sample delivered to LAB will be preceded or accompanied by complete written disclosure of the presence of any hazardous substances known or suspected by Client. Client further warrants that any sample containing any hazardous substance that is to be delivered to LAB will be packaged, labeled, transported, and delivered properly and in accordance with applicable laws.

Sample Handling.

Prior to LAB's acceptance of any sample (or after any revocation of acceptance), the entire risk of loss or of damage to such sample remains with Client. Samples are accepted when receipt is acknowledged on chain of custody documentation. In no event will LAB have any responsibility for the action or inaction of any carrier shipping or delivering any sample to or from LAB premises. Client authorizes LAB to proceed with the analysis of samples as received by the laboratory, recognizing that any samples not in compliance with all current DOH-ELAP-NELAP requirements for containers, preservation or holding time will be noted as such on the final report.

Disposal of hazardous waste samples is the responsibility of the Client. If the Client does not wish such samples returned, LAB may add storage and disposal fees to the final invoice. Maximum storage time for samples is 30 days after completion of analysis unless modified by applicable state or federal laws. Client will be required to give the LAB written instructions concerning disposal of these samples.

LAB reserves the absolute right, exercisable at any time, to refuse to receive delivery of, refuse to accept, or revoke acceptance of any sample, which, in the sole judgment of LAB (a) is of unsuitable volume, (b) may be or become unsuitable for or may pose a risk in handling, transport, or processing for any health, safety, environmental or other reason whether or not due to the presence in the sample of any hazardous substance, and whether or not such presence has been disclosed to LAB by Client or (c) if the condition or sample date make the sample unsuitable for analysis.

Legal Responsibility. LAB is solely responsible for performance of this contract, and no affiliated company, director, officer, employee, or agent shall have any legal responsibility hereunder, whether in contract or tort including negligence.

Assignment.

LAB may assign its performance obligations under this contract to other parties, as it deems necessary. LAB shall disclose to Client any assignee (subcontractor) by ELAP ID # on the submitted final report.

Force Majeure.

LAB shall have no responsibility or liability to the Client for any failure or delay in performance by LAB, which results in whole or in part from any cause or circumstance beyond the reasonable control of LAB. Such causes and circumstances shall include, but not limited to, acts of God, acts or orders of any government authority, strikes or other labor disputes, natural disasters, accidents, wars, civil disturbances, difficulties or delays in transportation, mail or delivery services, inability to obtain sufficient services or supplies from LAB's usual suppliers, or any other cause beyond LAB's reasonable control.

Law.

This contract shall be continued under the laws of the State of New York without regard to its conflicts of laws provision.

This report is part of a multipage document and should only be evaluated in its entirety. The Chain of Custody provides additional sample information, including compliance with the sample condition requirements upon receipt.

1066 Page 1 of 5 - KRM **CHAIN OF CUSTODY**

PARADIGM	CLIENT: NUI-VULL LL ADDRESS: 1667 LACE CITY COLLEGE STATE: N PHONE: (595) 470-1	CLIENT: ADDRESS: ZIP 146 15 CITY: BHONE: ATTN:	OICE TO: OTHER STATE: ZIP:	Quotation #:	yen-
ROPE GENERAL FMR. MGP SITE	Matrix Codes: AQ - Aqueous Liquid NQ - Non-Aqueous Liquid	WA - Water WG - Groundwater WW - Wastewater	er SL - Sludge	SD - Solid WP - Wipe PT - Paint CK - Caulk	OL - Oil AR - Air
DATE COLLECTED TIME COLLECTED COLLECTED S S I T E	SAMPLE IDENTIFIER	MATODES MAT		REMARKS	PARADIGM LAB SAMPLE NUMBER
723/22 14:00 × /23/22 12:45 × /23/22 - × 5/25/22 15:30 × /24/22 16:30 × /25/22 14:45 × /25/22 16:30 × /25/22 16	GEN-FB-0523 GEN-FB-0523 GEN-DUP-052 GEN-MWZ-05 GEN-MW 1-05 GEN-MW 9-05 GEN-MW 9-05 GEN-MW 9-05 GEN-MW 7-0 TIP Blank Report Supplements	2322 WG 3 XX 23 22 WG 3 XX 52322 WG 7 1 XX 52422 WG 3 XX 52422 WG 3 XX 2522 WG 3 XX 2522 WG 3 XX 2522 WG 3 XX	5/23-26/22	ignent Blank uplicate S/MSD	
andard 5 day day Batch QC Category A Category B Ish 1 day The Needed ase indicate date needed: None Required Category B Other please indicate packet	None Required Basic EDD NYSDEC EDD Other EDD	Received @ Lab By Custoday Sea Sintoct By signing this form, client agrees to	Date/Time Date/Time Signed, date d. Cf]

See additional page for sample conditions.

Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric **Park Street Former MGP Site**

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Parameter	Quantitation Limit ¹				
Volatile Organics	Water	Soil			
Method 8260	(µg/L)	(µg/kg) ²			
Chloromethane	5	5			
Bromomethane	5	5			
Vinyl Chloride	5	5			
Chloroethane	5	5			
Methylene Chloride	3	3			
Acetone	5	5			
Carbon Disulfide	5	5			
1,1-Dichloroethylene	5	5			
1,1-Dichloroethane	5	5			
1,2-Dichloroethylene (total)	5	5			
Chloroform	5	5			
1,2-Dichloroethane	2	2			
2-Butanone	5	5			
1,1,1-Trichloroethane	5	5			
Carbon Tetrachloride	2	2			
Bromodichloromethane	1	1			
1,2-Dichloropropane	1	1			
cis-1,3-Dichloropropene	5	5			
Trichloroethane	5	5			
Dibromochloromethane	5	5			
1,1,2-Trichloroethane	3	3			
Benzene	1	1			
trans-1,3-Dichloropropene	5	5			
Bromoform	4	4			
4-Methyl-2-pentanone	5	5			
2-Hexanone	5	5			
Tetrachloroethene	1	1			
Toluene	5	5			
1,1,2,2-Tetrachloroethane	1	1			
Chlorobenzene	5	5			
Ethylbenzene	4	4			
Styrene	5	5			
2-Chloroethyl Vinyl Ether	5	5			
1,2-Dichlorobenzene	5	5			
1,3-Dichlorobenzene	5	5			
1,4-Dichlorobenzene	5	5			
Vinyl Acetate	5	5			
Total Xylenes	5	5			
Semivolatile Organics	Water	Soil			
Method 8270	(µg/L)	(µg/kg)			
1,2,4-Trichlorobenzene	1	33			
1,2-Dichlorobenzene	10	330			
1,2-Diphenylhydrazine	10	330			
1,3-Dichlorobenzene	10	330			
1,4-Dichlorobenzene	10	330			
1,4-Dioxane	10	330			
2,4,5-Trichlorophenol	10	330			
2,4,6-Trichlorophenol	10	330			
2,4-Dichlorophenol	10	330			

Table 1 Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

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Parameter	Quantitation Limit ¹				
Semivolatile Organics	Water	Soil			
Method 8270 (Cont'd.)	(µg/L)	(µg/kg)			
2,4-Dimethylphenol	10	330			
2,4-Dinitrophenol	40	1300			
2,4-Dinitrotoluene	2	67			
2,6-Dinitrotoluene	2	67			
2-Chloronaphthalene	10	330			
2-Chlorophenol	10	330			
2-Methylnaphthalene	10	330			
2-Methylphenol	10	330			
2-Nitroaniline	20	670			
2-Nitrophenol	10	330			
3,3'-Dichlorobenzidene	20	670			
3-Nitroaniline	20	670			
4,6-Dinitro-2-methylphenol	40	1300			
4-Bromophenyl-phenylether	10	330			
4-Chioro-3-methylphenol	10	330			
4-Chloroaniline	10	330			
4-Chlorophenyl-phenylether	10	330			
4-Methylphenol	10	330			
4-Nitroaniline	20	670			
4-Nitrophenol	40	1300			
Acenaphthene	10	330			
Acenaphthylene	10	330			
Acetophenone	10	330			
Aniline	10	330			
Anthracene	10	330			
Atrazine	10	330			
Benzaldehyde	10	330			
Benzidine	40	1300			
Benzo(a)anthracene	1	33			
Benzo(a)pyrene	1	33			
Benzo(b)fluoranthene	1	33			
Benzo(g,h,i)perylene	10	330			
Benzo(k)fluoranthene	1	33			
Benzoic Acid	10	330			
Benzyl Alcohol	10	330			
bis(2-chloroethoxy)methane	10	330			
bis(2-chloroethyl)ether	1	33			
bis(2-chloroisopropyl)ether	10	330			
bis(2-ethylhexyl)phthalate	10	330			
Butylbenzylphthalate	10	330			
Caprolactam	10	330			
Carbazole	10	330			
Chrysene	10	330			
Dibenzo(a,h)anthracene	1	33			
Dibenzofuran	10	330			
Diethylphthalate	10	330			
Dimethylphthalate	10	330			
Di-n-butyl phthalate	10	330			
Di-n-octyl phthalate	10	330			

Table 1
Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

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Parameter	Quantitation Limit ¹			
Semivolatiles Method 8270 (Cont'd.)	Water (μg/L)	Soil (µg/kg)		
Diphenyl	10	330		
Fluoranthene	10	330		
Fluorene	10	330		
Hexachlorobenzene	1	33		
Hexachlorobutadiene	2	67		
Hexachlorocyclopentadiene	10	330		
Hexachloroethane	1	33		
Indeno(1,2,3-cd)pyrene	1	33		
Isophorone	10	330		
N,N-Dimethylaniline	1	33		
Naphthalene	10	330		
Nitrobenzene	1	33		
N-Nitrosodimethylamine	10	330		
N-Nitroso-di-n-propylamine	1	33		
N-Nitrosodiphenylamine	10	330		
Pentachlorophenol	40	1300		
Phenanthrene	10	330		
Phenol	10	330		
Pyrene	10	330		
Pyridine	10	330		
TAL Metals (6010/7470)	Water (µg/L)	Soil (µg/kg)		
Aluminum	(µg/L)	40		
Antimony		2		
Arsenic		1		
Barium	-	40		
Beryllium		0.4		
Cadmium		1		
Calcium	R=	1000		
Chromium		2		
Cobalt	***	10		
Copper		5		
Iron	76	30		
Lead		1		
Magnesium	W.F.	1000		
Manganese		3		
Mercury		0.033		
Nickel		8		
Potassium		1000		
Selenium		1000		
	-	2		
Silver Sodium				
		1000		
Thallium		2		
Vanadium		10		
Zinc	107	6		
Supplemental Parameters	Water (µg/L)	Soil (mg/kg)		
Total Organic Carbon (Lloyd Kahn)	NA	100		
Chloride Method 325.3	1,000			

Table 1 V
Parameter, Methods, and Quantitation Limits

Generic Quality Assurance Project Plan Rochester Gas & Electric Park Street Former MGP Site

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Parameter	Quantitation Limit ¹			
Supplemental Parameters (Cont'd.)	Water (μg/L)	Soil (mg/kg)		
Nitrate Method 353.2	100			
Ammonia Method 350.1	100			
Iron Method 200.7	150			
Manganese Method 200.7	15			
Sulfate Method 375.4	5,000			
Sulfide Method 376.1	1,000			
Orthophosphate Method 365.2	30			
Alkalinity Method 310.1	5,000			
Methane Method 3810	-			
Reactive Sulfide		20		
Reactive Cyanide		25		
TCLP Benzene		1		
Total Sulfur		50		
Chemical Oxygen Demand		120		

Notes:

¹ Specific quantitation limits are highly matrix dependent. The quantitation limits listed are for guidance and may not always be achievable due to matrix interference.

µg/L = micrograms per liter µg/kg = micrograms per kilogram

mg/kg = milligrams per kilogram

²Quantitation limits for source materials/soil/sediment are based on wet weight. The quantitation limits calculated by the laboratory for source materials/soil/sediment (calculated on a dry weight basis) will be higher.



Chain of Custody Supplement

Client:	Neu-Verre	Completed by:	Emilee Hyde
Lab Project ID:	222457	Date:	92012
	Sample Conditio Per NELAC/ELAP 21	on Requirements 0/241/242/243/244	
Condition	NELAC compliance with the sample of Yes	condition requirements t No	upon receipt N/A
Container Type Comments			
Transferred to method- compliant container			
Headspace (<1 mL) Comments	VOA		
Preservation Comments	VOA		
Chlorine Absent <0.10 ppm per test strip) Comments			
lolding Time Comments	4		
emperature Comments	D°C iced		
ompliant Sample Quantity/Ty Comments	уре		



ANALYTICAL REPORT

Lab Number: L2228058

Client: Paradigm Environmental Services

179 Lake Avenue Rochester, NY 14608

ATTN: Jane Daloia Phone: (585) 647-2530

Project Name: 222457
Project Number: 222457
Report Date: 06/09/22

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Lab Number: Report Date:

L2228058 06/09/22

Project Name: 222457
Project Number: 222457

05/26/22	05/26/22 13:00	Not Specified	WATER	GEN-MW7-052622 222457-09	L2228058-09
05/26/22	05/25/22 16:30	Not Specified	WATER	GEN-MW8-052522 222457-08	L2228058-08
05/26/22	05/25/22 14:45	Not Specified	WATER	GEN-MW6-052522 222457-07	L2228058-07
05/26/22	05/24/22 16:30	Not Specified	WATER	GEN-MW4-052422 222457-06	L2228058-06
05/26/22	05/24/22 14:45	Not Specified	WATER	GEN-MW1-052422 222457-05	L2228058-05
05/26/22	05/23/22 15:30	Not Specified	WATER	GEN-MW2-052322 222457-04	L2228058-04
05/26/22	05/23/22 00:00	Not Specified	WATER	GEN-DUP-052322 222457-03	L2228058-03
05/26/22	05/23/22 12:45	Not Specified	WATER	GEN-EB-052322 222457-02	L2228058-02
05/26/22	05/23/22 14:00	Not Specified	WATER	GEN MW3-052322 222457-01	L2228058-01
Receive Date	Collection Date/Time	Sample Location	Matrix	Client ID	Alpha Sample ID



 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Cattlin Wallet Caitlin Walukevich

Authorized Signature:

Title: Technical Director/Representative

Date: 06/09/22



ORGANICS



SEMIVOLATILES



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-01 Date Collected: 05/23/22 14:00

Client ID: GEN MW3-052322 222457-01 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/27/22 03:43
Analytical Date: 05/29/22 16:19

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM - Westborough Lab								
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptal Qualifier Criteri	
Nitrobenzene-d5	98	23-1	20
2-Fluorobiphenyl	78	15-1	20
4-Terphenyl-d14	70	41-1	49



Project Name: Lab Number: 222457 L2228058

Project Number: Report Date: 222457 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-02 Date Collected: 05/23/22 12:45

Date Received: 05/26/22 Client ID: GEN-EB-052322 222457-02 Sample Location: Field Prep: Not Specified Not Specified

05/29/22 16:35

Sample Depth:

Extraction Method: EPA 3510C Matrix: Water

Extraction Date: 05/27/22 03:43 Analytical Method: 1,8270D-SIM Analytical Date:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM	Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	114	23-120	
2-Fluorobiphenyl	87	15-120	
4-Terphenyl-d14	93	41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-03 Date Collected: 05/23/22 00:00

Client ID: GEN-DUP-052322 222457-03 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/27/22 03:43
Analytical Date: 05/29/22 16:52

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM	Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	88	23-120	
2-Fluorobiphenyl	70	15-120	
4-Terphenyl-d14	66	41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-04 Date Collected: 05/23/22 15:30

Client ID: GEN-MW2-052322 222457-04 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/27/22 03:43
Analytical Date: 05/29/22 17:08

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-S	Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	 1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	88	23-120	
2-Fluorobiphenyl	71	15-120	
4-Terphenyl-d14	66	41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-05 Date Collected: 05/24/22 14:45

Client ID: GEN-MW1-052422 222457-05 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/27/22 08:09
Analytical Date: 05/29/22 18:46

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM	Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	ND		ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	83	23-120	
2-Fluorobiphenyl	69	15-120	
4-Terphenyl-d14	72	41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-06 Date Collected: 05/24/22 16:30

Client ID: GEN-MW4-052422 222457-06 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/27/22 08:09
Analytical Date: 05/29/22 19:02

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor		
Semivolatile Organics by GC/MS-SIM	Semivolatile Organics by GC/MS-SIM - Westborough Lab							
Acenaphthene	ND		ug/l	0.10	0.04	1		
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1		
Fluoranthene	ND		ug/l	0.10	0.04	1		
Naphthalene	ND		ug/l	0.10	0.04	1		
Benzo(a)anthracene	ND		ug/l	0.10	0.02	1		
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1		
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1		
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1		
Chrysene	ND		ug/l	0.10	0.04	1		
Acenaphthylene	ND		ug/l	0.10	0.04	1		
Anthracene	ND		ug/l	0.10	0.04	1		
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1		
Fluorene	ND		ug/l	0.10	0.04	1		
Phenanthrene	0.03	J	ug/l	0.10	0.02	1		
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1		
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1		
Pyrene	ND		ug/l	0.10	0.04	1		
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
Nitrobenzene-d5	110	23-120	
2-Fluorobiphenyl	88	15-120	
4-Terphenyl-d14	90	41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-07 Date Collected: 05/25/22 14:45

Client ID: GEN-MW6-052522 222457-07 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/29/22 11:36
Analytical Date: 05/31/22 20:01

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Semivolatile Organics by GC/MS-SIM	- Westborough La	ab					
Acenaphthene	2.9		ug/l	0.10	0.04	1	
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1	
Fluoranthene	0.05	J	ug/l	0.10	0.04	1	
Naphthalene	200	Е	ug/l	0.10	0.04	1	
Benzo(a)anthracene	0.02	J	ug/l	0.10	0.02	1	
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1	
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1	
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1	
Chrysene	ND		ug/l	0.10	0.04	1	
Acenaphthylene	29	Е	ug/l	0.10	0.04	1	
Anthracene	0.42		ug/l	0.10	0.04	1	
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1	
Fluorene	5.5		ug/l	0.10	0.04	1	
Phenanthrene	1.9		ug/l	0.10	0.02	1	
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1	
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1	
Pyrene	ND		ug/l	0.10	0.04	1	
2-Methylnaphthalene	0.10	J	ug/l	0.10	0.05	1	

Surrogate	% Recovery	Qualifier	Acceptance Criteria	
Nitrobenzene-d5	110		23-120	
2-Fluorobiphenyl	80		15-120	
4-Terphenyl-d14	70		41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-07 D Date Collected: 05/25/22 14:45

Client ID: GEN-MW6-052522 222457-07 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/29/22 11:36
Analytical Date: 06/09/22 16:56

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM - West	oorough Lab					
Naphthalene	180		ug/l	1.0	0.43	10
Acenaphthylene	21		ug/l	1.0	0.35	10



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-08 Date Collected: 05/25/22 16:30

Client ID: GEN-MW8-052522 222457-08 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/29/22 11:36
Analytical Date: 05/31/22 20:17

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-SIM	- Westborough La	b				
Acenaphthene	ND		ug/l	0.10	0.04	1
2-Chloronaphthalene	ND		ug/l	0.20	0.04	1
Fluoranthene	ND		ug/l	0.10	0.04	1
Naphthalene	ND		ug/l	0.10	0.04	1
Benzo(a)anthracene	0.02	J	ug/l	0.10	0.02	1
Benzo(a)pyrene	ND		ug/l	0.10	0.04	1
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04	1
Chrysene	ND		ug/l	0.10	0.04	1
Acenaphthylene	ND		ug/l	0.10	0.04	1
Anthracene	ND		ug/l	0.10	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.10	0.04	1
Fluorene	ND		ug/l	0.10	0.04	1
Phenanthrene	ND		ug/l	0.10	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04	1
Pyrene	ND		ug/l	0.10	0.04	1
2-Methylnaphthalene	ND		ug/l	0.10	0.05	1

Surrogate	% Recovery	Qualifier	Acceptance Criteria	
Nitrobenzene-d5	90		23-120	
2-Fluorobiphenyl	62		15-120	
4-Terphenyl-d14	62		41-149	



Project Name: 222457 Lab Number: L2228058

Project Number: 222457 Report Date: 06/09/22

SAMPLE RESULTS

Lab ID: L2228058-09 Date Collected: 05/26/22 13:00

Client ID: GEN-MW7-052622 222457-09 Date Received: 05/26/22 Sample Location: Not Specified Field Prep: Not Specified

Sample Depth:

Matrix: Water Extraction Method: EPA 3510C

Analytical Method: 1,8270D-SIM Extraction Date: 05/29/22 11:36
Analytical Date: 05/31/22 20:34

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Semivolatile Organics by GC/MS-S	IM - Westborough La	b				
Acenaphthene	ND		ug/l	0.11	0.04	1
2-Chloronaphthalene	ND		ug/l	0.22	0.04	1
Fluoranthene	ND		ug/l	0.11	0.04	1
Naphthalene	ND		ug/l	0.11	0.05	1
Benzo(a)anthracene	ND		ug/l	0.11	0.02	1
Benzo(a)pyrene	ND		ug/l	0.11	0.04	1
Benzo(b)fluoranthene	ND		ug/l	0.11	0.02	1
Benzo(k)fluoranthene	ND		ug/l	0.11	0.05	1
Chrysene	ND		ug/l	0.11	0.04	1
Acenaphthylene	ND		ug/l	0.11	0.04	1
Anthracene	ND		ug/l	0.11	0.04	1
Benzo(ghi)perylene	ND		ug/l	0.11	0.05	1
Fluorene	ND		ug/l	0.11	0.04	1
Phenanthrene	0.03	J	ug/l	0.11	0.02	1
Dibenzo(a,h)anthracene	ND		ug/l	0.11	0.04	1
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.11	0.04	1
Pyrene	ND		ug/l	0.11	0.04	1
2-Methylnaphthalene	ND		ug/l	0.11	0.05	1

Surrogate	% Recovery	Qualifier	Acceptance Criteria	
Nitrobenzene-d5	116		23-120	
2-Fluorobiphenyl	77		15-120	
4-Terphenyl-d14	71		41-149	



Project Name: 222457

Project Number: 222457

Lab Number:

L2228058

Report Date: 06/09/22

Method Blank Analysis Batch Quality Control

Analytical Method: Analytical Date:

1,8270D-SIM 05/29/22 16:03

Analyst:

RP

Extraction Method: EPA 3510C Extraction Date: 05/27/22 03:43

arameter	Result	Qualifier Units	RL	MDL	
emivolatile Organics by GC/M	S-SIM - Westbo	rough Lab for samp	le(s): 01-04	Batch: WG1	643559-
Acenaphthene	ND	ug/l	0.10	0.04	
2-Chloronaphthalene	ND	ug/l	0.20	0.04	
Fluoranthene	ND	ug/l	0.10	0.04	
Naphthalene	ND	ug/l	0.10	0.04	
Benzo(a)anthracene	ND	ug/l	0.10	0.02	
Benzo(a)pyrene	ND	ug/l	0.10	0.04	
Benzo(b)fluoranthene	ND	ug/l	0.10	0.02	
Benzo(k)fluoranthene	ND	ug/l	0.10	0.04	
Chrysene	ND	ug/l	0.10	0.04	
Acenaphthylene	ND	ug/l	0.10	0.04	
Anthracene	ND	ug/l	0.10	0.04	
Benzo(ghi)perylene	ND	ug/l	0.10	0.04	
Fluorene	ND	ug/l	0.10	0.04	
Phenanthrene	ND	ug/l	0.10	0.02	
Dibenzo(a,h)anthracene	ND	ug/l	0.10	0.04	
Indeno(1,2,3-cd)pyrene	ND	ug/l	0.10	0.04	
Pyrene	ND	ug/l	0.10	0.04	
2-Methylnaphthalene	ND	ug/l	0.10	0.05	

		Acceptance
Surrogate	%Recovery	Qualifier Criteria
Nitrobenzene-d5	91	23-120
2-Fluorobiphenyl	71	15-120
4-Terphenyl-d14	87	41-149



Project Name: 222457

Project Number: 222457

Lab Number:

L2228058

Report Date: 06/09/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D-SIM Analytical Date: 05/29/22 18:30

Analyst:

RP

Extraction Method: EPA 3510C Extraction Date: 05/27/22 08:09

Qualifier RL MDL **Parameter** Result Units Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 05-06 Batch: WG1643644-1 Acenaphthene ND ug/l 0.10 0.04 2-Chloronaphthalene ND ug/l 0.20 0.04 ND 0.10 0.04 Fluoranthene ug/l Naphthalene ND ug/l 0.10 0.04 ND 0.02 Benzo(a)anthracene ug/l 0.10 ND 0.04 Benzo(a)pyrene ug/l 0.10 Benzo(b)fluoranthene ND ug/l 0.10 0.02 Benzo(k)fluoranthene ND ug/l 0.10 0.04 Chrysene ND 0.10 0.04 ug/l ND 0.04 Acenaphthylene 0.10 ug/l 0.10 0.04 Anthracene ND ug/l Benzo(ghi)perylene ND ug/l 0.10 0.04 Fluorene ND ug/l 0.10 0.04 Phenanthrene ND ug/l 0.10 0.02 Dibenzo(a,h)anthracene ND ug/l 0.04 0.10 ND Indeno(1,2,3-cd)pyrene ug/l 0.10 0.04

ug/l

ug/l

0.10

0.10

		Acceptance
Surrogate	%Recovery	Qualifier Criteria
Nitrobenzene-d5	79	23-120
2-Fluorobiphenyl	65	15-120
4-Terphenyl-d14	72	41-149

ND

ND



0.04

0.05

Pyrene

2-Methylnaphthalene

Project Name: 222457

Project Number: 222457

Lab Number:

L2228058

Report Date: 06/09/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D-SIM Analytical Date: 05/31/22 19:45

Analyst:

RP

Extraction Method: EPA 3510C Extraction Date: 05/29/22 11:36

arameter	Result	Qualifier	Units	RL	MDL
emivolatile Organics by GC/MS-	SIM - Westbo	rough Lab	for sample(s)	: 07-09	Batch: WG1644257-1
Acenaphthene	ND		ug/l	0.10	0.04
2-Chloronaphthalene	ND		ug/l	0.20	0.04
Fluoranthene	ND		ug/l	0.10	0.04
Naphthalene	ND		ug/l	0.10	0.04
Benzo(a)anthracene	0.02	J	ug/l	0.10	0.02
Benzo(a)pyrene	ND		ug/l	0.10	0.04
Benzo(b)fluoranthene	ND		ug/l	0.10	0.02
Benzo(k)fluoranthene	ND		ug/l	0.10	0.04
Chrysene	ND		ug/l	0.10	0.04
Acenaphthylene	ND		ug/l	0.10	0.04
Anthracene	ND		ug/l	0.10	0.04
Benzo(ghi)perylene	ND		ug/l	0.10	0.04
Fluorene	ND		ug/l	0.10	0.04
Phenanthrene	ND		ug/l	0.10	0.02
Dibenzo(a,h)anthracene	ND		ug/l	0.10	0.04
Indeno(1,2,3-cd)pyrene	ND		ug/l	0.10	0.04
Pyrene	ND		ug/l	0.10	0.04
2-Methylnaphthalene	ND		ug/l	0.10	0.05

		Acceptance
Surrogate	%Recovery	Qualifier Criteria
Nitrobenzene-d5	108	23-120
2-Fluorobiphenyl	73	15-120
4-Terphenyl-d14	77	41-149



Project Number: 222457

Project Name:

222457

Lab Number: L2228058

Parameter	%Recovery	Qual	%Recovery	%Recovery Qual Limits	ry RPD	Qual	Limits
Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG1643559-2 WG1643559-3	stborough Lab As	sociated sam	ple(s): 01-04	Batch: WG1643559-2	WG1643559-3		
Acenaphthene	72		83	40-140	14		40
2-Chloronaphthalene	76		86	40-140	12		40
Fluoranthene	84		94	40-140	11		40
Naphthalene	70		82	40-140	16		40
Benzo(a)anthracene	85		96	40-140	12		40
Benzo(a)pyrene	100		112	40-140	11		40
Benzo(b)fluoranthene	89		94	40-140	5		40
Benzo(k)fluoranthene	89		104	40-140	16		40
Chrysene	75		86	40-140	14		40
Acenaphthylene	91		103	40-140	12		40
Anthracene	82		93	40-140	13		40
Benzo(ghi)perylene	81		90	40-140	11		40
Fluorene	80		91	40-140	13		40
Phenanthrene	74		83	40-140	11		40
Dibenzo(a,h)anthracene	87		97	40-140	11		40
Indeno(1,2,3-cd)pyrene	82		92	40-140	11		40
Pyrene	82		92	40-140	1		40
2-Methylnaphthalene	75		86	40-140	14		40



Project Number: 222457

Project Name:

222457

Lab Number: L2228058

	LCS		LCSD		%Recovery			RPD	
Parameter	%Recovery Qual	Qual	%Recovery	Qual	Limits	RPD	Qual	Limits	
Semivolatile Organics by GC/MS-SIM - Wes	sthorough I ah	Associated san	nnle/s): 01-04	Ratch: 1	3550-0	NG1643550-3			
Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG164:	stborough Lab	Associated san	nple(s): 01-04	Batch: \	3559-2	WG1643559-3			

Nitrobenzene-d5 2-Fluorobiphenyl 4-Terphenyl-d14	Surrogate
105 74 79	LCS %Recovery Qual
124 85 88	LCSD Qual %Recovery
Q	y Qual
23-120 15-120 41-149	Acceptance Criteria



Project Name: 222457

Project Number: 222457

> Lab Number: L2228058

	LCS %Bccycer	2	LCSD %Recovery	2	%Recovery	B B D	2	RPD	
Parameter	%Recovery	%Recovery Qual	%Recovery Qual	Qual	Limits	RPD	Qual	Limits	
Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 05-06 Batch: WG164	Vestborough Lab	Associated samp	le(s): 05-06	Batch: Wo	31643644-2 W	43644-2 WG1643644-3			

			,	****		:	4	
semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 05-06 Batch: WG1643644-2 WG1643644-3	orough Lab ,	Associated sample(s): 05-06	Batch: WG1643644-	2 WG1643	8644-3		
Acenaphthene	84		64	40-140		27		40
2-Chloronaphthalene	82		64	40-140		25		40
Fluoranthene	93		72	40-140		25		40
Naphthalene	82		64	40-140	0	25		40
Benzo(a)anthracene	87		66	40-140	0	27		40
Benzo(a)pyrene	109		84	40-140	0	26		40
Benzo(b)fluoranthene	92		70	40-140	0	27		40
Benzo(k)fluoranthene	94		73	40-140	0	25		40
Chrysene	87		68	40-140	0	25		40
Acenaphthylene	96		74	40-140		26		40
Anthracene	90		69	40-140		26		40
Benzo(ghi)perylene	98		73	40-140		29		40
Fluorene	88		67	40-140		27		40
Phenanthrene	81		62	40-140		27		40
Dibenzo(a,h)anthracene	103		78	40-140		28		40
Indeno(1,2,3-cd)pyrene	96		74	40-140		26		40
Pyrene	93		72	40-140		25		40
2-Methylnaphthalene	85		66	40-140	0	25		40



Lab Number:

L2228058

Project Number: 222457 Project Name:

222457

Report Date: 06/09/22

Parameter	LCS %Recovery Qual	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Combinatella Dissociate has OCMO SIM Wheelt-course like Approjected completely of OC Details WO424				Dotob: 141		0.000.00			

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 05-06 Batch: WG1643644-2 WG1643644-3

Nitrobenzene-d5 2-Fluorobiphenyl 4-Terphenyl-d14	Surrogate
114 83 89	LCS %Recovery Qual
90 65 71	LCSD %Recovery
23-120 15-120 41-149	Acceptance Qual Criteria



Project Number: Project Name:

222457 222457

> Lab Number: L2228058

Parameter	LCS %Recovery	Qual	LCSD %Recovery	%Recovery Qual Limits	ery RPD	Qual	RPD Limits
Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 07-09 Batch: WG1644257-2 WG1644257-3	tborough Lab As	ssociated sam	ple(s): 07-09	Batch: WG1644257-2	WG1644257-3		
Acenaphthene	92		75	40-140	20		40
2-Chloronaphthalene	96		78	40-140	21		40
Fluoranthene	103		86	40-140	18		40
Naphthalene	94		76	40-140	21		40
Benzo(a)anthracene	110		93	40-140	17		40
Benzo(a)pyrene	123		101	40-140	20		40
Benzo(b)fluoranthene	111		90	40-140	21		40
Benzo(k)fluoranthene	98		82	40-140	18		40
Chrysene	93		78	40-140	18		40
Acenaphthylene	109		89	40-140	20		40
Anthracene	101		83	40-140	20		40
Benzo(ghi)perylene	113		95	40-140	17		40
Fluorene	99		81	40-140	20		40
Phenanthrene	94		76	40-140	21		40
Dibenzo(a,h)anthracene	121		100	40-140	19		40
Indeno(1,2,3-cd)pyrene	133		110	40-140	19		40
Pyrene	99		83	40-140	18		40
2-Methylnaphthalene	98		79	40-140	21		40



Project Number: 222457

Project Name:

222457 Report Date: Lab Number: 06/09/22 L2228058

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 07-09 Batch: WG164	Parameter
IM - Westborough Lab As	LCS %Recovery
ssociated sa	Qual
ample(s): 07-0	LCSD %Recover
9 Batch:	y Qua
WG1644257-2	%Recove
4257-2 WG1644257-3	yry RPD
	Qual
	RPD Limits

	LCS		LCSD		Acceptance
Surrogate	%Recovery Qual	Qual	%Recovery Qual	Qual	Criteria
Nitrobenzene-d5	142	Q	111		23-120
2-Fluorobiphenyl	94		75		15-120
4-Terphenyl-d14	96		78		41-149



Matrix Spike Analysis Batch Quality Control

Project Number: Project Name: 222457 222457

Lab Number:

Report Date:

06/09/22 L2228058

058-04	ole: L2228	QC Samp	43559-5	WG16	WG1643559-4	QC Batch ID:	1-04	ated sample(s): 0	Associa	stborough Lab	Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 QC Batch ID: WG1643559-4 WG1643559-5 QC Sample: L2228058-04	Semivolatil
RPD Limits	Qual L	y RPD	Recovery Limits	Qual	MSD Recovery RPD %Recovery Qual Limits RPD Qual Limits	MSD al Found	Qua	MS %Recovery Qual	MS Found	MS Added	Native Sample	Parameter

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Client ID: GEN-MW2-052322 222457-04	Л - Westborou 7-04	gh Lab A	ssociated sa	mple(s): 01-04	QC Batch ID: W	ı ID: WG1	643559-4	WG1643	559-5 QC	Sample: L	'G1643559-4 WG1643559-5 QC Sample: L2228058-04
Acenaphthene	ND 10	0	6.4	64	7.8	8	78		40-140	20	40
2-Chloronaphthalene	ND 10	0	6.3	63	7.7	7	77		40-140	20	40
Fluoranthene	ND 10	0	7.8	78	9.3	ω	93		40-140	18	40
Naphthalene	ND 10	0	6.2	62	7.5	5	75		40-140	19	40
Benzo(a)anthracene	ND 10	0	7.0	70	8.5	ST.	85		40-140	19	40
Benzo(a)pyrene	ND 10	0	4.4	44	5.0	0	50		40-140	13	40
Benzo(b)fluoranthene	ND 10	0	4.2	42	5.0	0	50		40-140	17	40
Benzo(k)fluoranthene	ND 10	0	3.8	38	Q 4.2	2	42		40-140	10	40
Chrysene	ND 10	0	5.3	53	6.2	2	62		40-140	16	40
Acenaphthylene	ND 10	0	6.9	69	8.4	4	84		40-140	20	40
Anthracene	ND 10	0	6.7	67	8.1		81		40-140	19	40
Benzo(ghi)perylene	ND 10	0	1.2	12	Q 1.4	4	14	۵	40-140	15	40
Fluorene	ND 10	0	6.8	68	8.2	2	82		40-140	19	40
Phenanthrene	ND 10	0	6.1	61	7.5	5	75		40-140	21	40
Dibenzo(a,h)anthracene	ND 10	0	1.4	14	Q 1.6	6	16	۵	40-140	13	40
Indeno(1,2,3-cd)pyrene	ND 10	0	1.4	14	Q 1.7	7	17	۵	40-140	19	40
Pyrene	ND 10	0	8.1	81	9.5	5	95		40-140	16	40
2-Methylnaphthalene	ND 10	0	6.2	62	7.5	5	75		40-140	19	40

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
2-Fluorobiphenyl	60	73	15-120
4-Terphenyl-d14	58	64	41-149



Matrix Spike Analysis

Batch Quality Control

Report Date: Lab Number:

06/09/22 L2228058

Project Number: 222457

Project Name:

222457

Parameter Native Sample MS Added MS Found %Recovery Qual MSD MSD Found %Recovery Qual Recovery Limits RPD Qual Limits

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 QC Batch ID: WG1643559-4 WG1643559-5 QC Sample: L2228058-04 Client ID: GEN-MW2-052322 222457-04

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
Nitrobenzene-d5	68	84	23-120



Project Name: 222457
Project Number: 222457

Report Date: 06/09/22

Serial_No:06092219:02 **Lab Number:** L2228058

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal

Absent

Container Information Container ID Contai	rmation Container Type	Cooler	Initial pH	Final pH	Temp deg C Pres Seal	Pres	Seal	Frozen Date/Time	Analysis(*)
L2228058-01A	Amber 1000ml unpreserved	A	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-02A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-03A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-04A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-04A1	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-04A2	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-05A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-06A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-07A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-08A	Amber 1000ml unpreserved	≻	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)
L2228058-09A	Amber 1000ml unpreserved	Þ	7	7	3.8	~	Absent		NYTCL-PAHSIM(7)



 Project Name:
 222457
 Lab Number:
 L2228058

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GLOSSARY

Acronyms

EDL

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

 Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

EMPC - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content,

where applicable. (DoD report formats only.)

LOQ - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

only.)

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any

adjustments from dilutions, concentrations or moisture content, where applicable.

MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

NR - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile

Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEQ - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF

and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Footnotes

1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

- A -Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Data Qualifiers

- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- V The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
 (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Serial_No:06092219:02

 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Serial_No:06092219:02

Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

Revision 19 Published Date: 4/2/2021 1:14:23 PM

ID No.:17873

Page 1 of 1

Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

EPA 8260C/8260D: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522, EPA 537.1.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Pre-Qualtrax Document ID: 08-113 Document Type: Form

179 Lake Avenue, Rochester, NY 14608 Office (585) 647-2530 Fax (585) 647-3311

CHAIN OF CUSTODY L2228058

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1 5/23/22	14:00		X	GEN-MW3-05	2322	WG	1	x	NYS	ASP Cat B	222	457-01			
2 5/23/22	12:45		Х	GEN-EB-0523	322	WG	1	X	Rep	ort Pkg and		02			
3 5/23/22			X	GEN-DUP-05	2322	WG	3	x	NYS	SDEC EDD		03			
4 5/23/22	15:30		X	GEN-MW2-05	2322	WG	1	x	SIM fo	r PAH DLs as		64			
5 5/24/22	14:45		X	GEN-MW1-05	2422	WG	1	x		needed		65			
6 5/24/22	16:30		Х	GEN-MW4-05	2422	WG	4	x				06			
7 5/25/22	14:45		х	GEN-MW6-05	2522	WG	্ৰ	x	Run	MS/MSD on		07			
8 5/25/22	16:30		X	GEN-MW8-05	2522	WG	1	x	141 1 A	e GEN-MW2-		08	,		
9 5/26/22	13:00		Х	GEN-MW7-05	2622	WG	1	x		052322		09			
10					04450										
The second second second second	on: Per NELAC/E Receipt Parame Container Type	LAP 210/24 iter	Children Co.	43/244 NELAC Comp		Client		E							
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Page 34 of 3	4				Rec	ceived @ Lab	Ву	_		Date/Time					

DATA USABILITY SUMMARY REPORT (DUSR)

RGE Geneseo Former MGP Site

SDGs: 222457

8 Water Samples, equipment blank and trip blank

Prepared for:

Neu Velle, LLC 1667 Lake Ave., Bldg. 59, Suite 101 Rochester, NY 14615 **Attention: Kyle Miller**

September 2022



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Tables

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Table 4-2 Quality Control Criteria for Validating Laboratory Analytical Data

Summaries of Validated Results

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REVIEWER'S NARRATIVE

Neu-Velle SDG 222457: RGE Geneseo Former MGP Site

The data associated with this Sample Delivery Group (SDG) 222457, analyzed by Paradigm Environmental Services, Inc. Rochester, NY have been reviewed in accordance with assessment criteria provided by the New York State Department of Environmental Conservation following the review procedures provided in the USEPA Functional Guidelines for evaluating organic and inorganic data.

All analytical results reported by the laboratory are considered valid and acceptable except results that have been qualified as rejected, "R". Results qualified as estimated "J", or as non-detects, "U", are considered usable for the purpose of evaluating water and/or soil quality. However, these qualifiers indicate that the accuracy and/or precision of the analytical result is questionable. A summary of all data that have been qualified and the reasons for qualification are provided in the following data usability summary report (DUSR).

Two facts should be noted by all data users. First, the "R" qualifier means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the analyte is present or not. Values qualified with an "R" should not appear on the final data tables because they cannot be relied upon, even as the last resort. Second, no analyte concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error.

Reviewer's Signature:	Míchael K. Perry	Date:	9/8/2022	
_	Michael K. Perry			
	Chemist			

1.0 SUMMARY

SITE: RGE Geneseo

Former MGP Site

SAMPLING DATE: May 23 – May 26, 2022

SAMPLE TYPE: 8 water samples, equipment and trip blank

LABORATORY: Paradigm Environmental Services, Inc.

Rochester, NY

SDG No.: 222457

2.0 INTRODUCTION

This data usability summary report (DUSR) was prepared in accordance with guidance provided by the New York State Department of Environmental Conservation (NYSDEC). The DUSR is based on a review and evaluation of the laboratory analytical data package. Specifically, the NYSDEC guidance recommends review and evaluation of the following elements of the data package:

- Completeness of the data package as defined under the requirements of the NYSDEC Analytical Services Protocols (ASP) Category B or the United States Environmental Protection Agency (USEPA) Contract Laboratory Program (CLP) deliverables,
- Compliance with established analyte holding times,
- Adherence to quality control (QC) limits and specifications for blanks, instrument tuning and calibration, surrogate recoveries, spike recoveries, laboratory duplicate analyses, and other QC criteria,
- Adherence to established analytical protocols,
- Conformance of data summary sheets with raw analytical data, and
- Use of correct data qualifiers.

Neu-Velle Page 1

Data deficiencies, analytical protocol deviations, and quality control problems identified using the review criteria above and their effect on the analytical results are discussed in this report.

3.0 SAMPLE AND ANALYSIS SUMMARY

The data packages consists of analytical results for eight water samples, equipment blank, and a trip blank collected on May 23 – May 26, 2022. These samples were analyzed for volatile organic compounds (BTEX) and Semi-Volatile Organic Compounds (SVOCs).

All analyses were performed by Paradigm Environmental Services, Inc., Rochester, NY and analyzed as SDG: 222457. The SVOCs were subcontracted to Alpha Analytical in Westborough, MA and analyzed as SDG: L2228058 for PAHs by 8270-D-SIM. The analytical results were provided in NYSDEC ASP Category B format, which includes all raw analytical data and laboratory QC data.

4.0 GUIDANCE DOCUMENTS AND DATA REVIEW CRITERIA

The guidance documents appropriate for reviewing laboratory quality control (QC) data and assigning data qualifiers (flags) to analytical results were selected from those listed in Table 4-1. The QC limits established in the documents applicable to this data review were used to assess the quality of the analytical results. In some cases, however, QC limits established internally by the laboratory were taken into account to determine data quality.

The QC criteria considered for assessing the usability of the reported analytical results provided for each analyte type (i.e. VOCs, SVOCs, metals, etc.) are listed in Table 4-2. These criteria may vary with the analytical method utilized by the laboratory. These criteria comply with the guidance recommended in Section 2.0 above.

5.0 DATA VALIDATION QUALIFIERS

The letter qualifiers (flags) used to define data usability are described briefly below. These letters are assigned by the data validator to analytical results having questionable accuracy and/or precision as determined by reviewing the laboratory QC data associated with the analytical results.

Neu-Velle Page 2

TABLE 4-1

Guidance Used For Validating Laboratory Analytical Data

Analyte Group	Guidance	Date
Metals (ICP-AES)	USEPA SOP HW-3a, Rev. 1	September 2016
Metals (Hg & CN)	USEPA SOP HW-3c, Rev. 1	September 2016
Volatile Organic Compounds (by Methods 8260B & 8260C)	USEPA SOP HW-24, Rev. 4	September 2014
Semi-Volatile Organic Compounds (by Method 8270D)	USEPA SOP HW-22 Rev. 5	December 2010
Pesticides (by Method 8181B)	USEPA SOP HW-44, Rev. 1.1	December 2010
Chlorinated Herbicides (by Method 8151A)	USEPA SOP HW-17, Rev. 3.1	December 2010
Polychlorinated Biphenyls (PCBs)	USEPA SOP HW-37A, Rev. 0	June 2015
Volatile Organic Compounds (Air) (by Method TO-15)	USEPA SOP HW-31, Rev. 6	September 2016
Per- and PolyFluoroAlkyl Substances (PFAS)	* NYSDEC	January 2021
General Chemistry Parameters	per NYSDEC ASP	July 2005

^{*} Sampling, Analysis, and Assessment of Per- and Polyfluoroalkyl Substances (PFAS) Under NYSDEC's Part 375 Remedial Programs, Appendix I

TABLE 4-2

QUALITY CONTROL CRITERIA USED FOR VALIDATING LABORATORY ANALYTICAL DATA

VOCs	SVOCs	Pesticides/PCBs	Metals	Gen Chemistry	PFAS
Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg	Completeness of Pkg
Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation	Sample Preservation
Holding Time	Holding Time	Holding Time	Holding Time	Holding Times	Holding Time
System Monitoring	Surrogate Recoveries	Surrogate Recoveries	Initial/Continuing	Calibration	Instr Performance
Compounds	Lab Control Sample	Matrix Spikes	Calibration	Lab Control Samples	Check
Lab Control Sample	Matrix Spikes	Blanks	CRDL Standards	Blanks	Initial Calibration
Matrix Spikes	Blanks	Instrument Calibration	Blanks	Spike Recoveries	Continuing Calibration
Blanks	Instrument Tuning	& Verification	Interference Check	Lab Duplicates	Blanks
Instrument Tuning	Internal Standards	Comparison of	Sample		Surrogates
Internal Standards	Initial Calibration	duplicate	Spike Recoveries		Lab Fortified Blank
Initial Calibration	Continuing Calibration	GC column results	Lab Duplicate		Matrix Spikes
Continuing Calibration	Lab Qualifiers	Analyte ID	Lab Control Sample		Internal Standards
Lab Qualifiers	Field Duplicate	Lab Qualifiers	ICP Serial Dilutions		
Field Duplicate		Field Duplicate	Lab Qualifiers		
			Field Duplicate		

Method TO-15 (Air)

Completeness of Pkg
Sample Preservation
Holding Time
Canister Certification
Instrument Tuning
Initial Calibration and
Instrument Performance
Daily Calibration
Blanks
Lab Control Sample
Field Duplicate

The laboratory may also use various letters and symbols to flag analytical results generated when QC limits were exceeded. The meanings of these flags may differ from those used by the independent data validator. Those used by the laboratory are provided with the analytical results.

NOTE: The assignment of data qualifiers by the data reviewer (validator) to laboratory analytical results should not necessarily be interpreted by the data user as a measure of laboratory ability or proficiency. Rather, the qualifiers are intended to provide a measure of data accuracy and precision to the data user, which, for example, may provide a level of confidence in determining whether or not standards or cleanup objectives have been met.

- U The analyte was analyzed for but was not detected at or above the sample quantitation limit.
- J The analyte was positively identified; the associated numerical value is the *approximate* concentration of the analyte in the sample. (The magnitude of any \pm value associated with the result is not determined by data validation).
- **J**+ The result is an estimated quantity and may be biased high.
- **J-** The result is an estimated quantity and may be biased low.
- **UJ** The analyte was analyzed for but not detected. The reported quantitation limit is approximate and may inaccurate or imprecise.
- R The sample result is rejected (i.e., is unusable) due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- **NJ** The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

The validated analytical results are attached to this report. Validation qualifiers (flags) are indicated in red print. Data sheets having qualified data are signed and dated by the data reviewer.

Neu-Velle Page 3

6.0 RESULTS OF THE DATA REVIEW

The results of the data review are summarized in Tables 6-1 through 6-2. The tables list the samples where QC criteria were found to exceed acceptable limits and the actions taken to qualify the associated analytical results.

7.0 TOTAL USABLE DATA

For SDG 222457, eight samples, equipment blank and a trip blank were analyzed and results were reported for 222 analytes. Even though some results were flagged with a "J" as estimated, all results (100%) are considered usable.

Neu-Velle Page 4

Table 6-1 8260 - BTEX

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
GEN-EB-5/23/22 GEN-MW6-5/25/22 GEN-MW8-5/25/22 Trip Blank	All analytes	J detects UJ non-detects	Surr rec for 1,2-DCEd4 < QC limit	Data are estimated
GEN-DUP-5/23/22 GEN-MW2-5/23/22 GEN-MW1-5/24/22 GEN-MW4-5/25/22	All analytes	J detects UJ non-detects	Surr rec for 1,2-DCEd4 and Td8 < QC limit	Data are estimated
GEN-MW2-5/23/22	Ethyl benzene	J detects	MS/MSD > QC limit	Detected data are estimated

Table 6-2 8270-SIM-PAH

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
GEN-MW6-5/25/22 GEN-MW8-5/25/22	Benzo(a)anthracene	CRQL-U	Detected in method blank	Detected data changed to non-detect
GEN-MW2-5/23/22	Benzo(k)fluoranthene Benzo(g,h,i)perylene Dibenzo(a,h)anthracene Indene(1,2,3-cd)pyrene	J detects UJ non-detects	MS/MSD < QC limit	Data are estimated
GEN-MW3-5/23/22 GEN-EB-5/23/22 GEN-DUP-5/23/22 GEN-MW2-5/23/22 GEN-MW1-5/24/22 GEN-MW4-5/25/22	Benzo(a)pyrene	J detects UJ non-detects	CCV < QC limit	Data are estimated

SDG 222457

GEN-MW6-5/25/22 GEN-MW8-5/25/22 GEN-MW7-5/26/22 Benzo(a)pyrene Indene(1,2,3-cd)pyrene	J detects UJ non-detects	CCV < QC limit	Data are estimated
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ACRONYMS

BSP

Blank Spike

CCAL

Continuing Calibration

CCB

Continuing Calibration Blank

CCV

Continuing Calibration Verification

CRDL

Contract Required Detection Limit

CRQL

Contract Required Quantitation Limit

% D

Percent Difference

ICAL

Initial Calibration

ICB

Initial Calibration Blank

IS

Internal Standard

LCS

Laboratory Control Sample

MS/MSD

Matrix Spike/Matrix Spike Duplicate

QA

Quality Assurance

QC

Quality Control

%R

Percent recovery

RPD

Relative Percent Difference

RRF

Relative Response Factor

%RSD

Percent Relative Standard Deviation

TAL

Target Analyte List (metals)

TCL

Target Compound List (organics)

Appendix A

Validated Analytical Results

LAB PROJECT NARRATIVE: 222457 PROJECT NAME: RGE Geneseo Fmr. MGP Site

SDG: 2457-01 CLIENT: Neu-Velle

Seven groundwater samples were collected by the client on May 23, 2022 and were received by the Paradigm laboratory on May 26, 2022. Samples were accompanied by a field duplicate, equipment blank, and trip blank. The samples were received under the conditions as noted on the chain-of-custody supplement. The samples were submitted with the Chains-of-Custody requesting the BTEX list for Volatiles and PAH by 8270SIM. All analyses, where applicable, were performed using EPA SW-846 Methods and the associated holding times.

The items noted in this case narrative address compliance with the referenced methods, NYSDOH ELAP rules, and any project specific data quality requirements. These may be different from the usability criteria referenced in any "Functional Guidelines" or other data review standards used by data validators.

GENERAL NOTES

ALL ANALYSES

The initial and continuing calibration reports are only evaluated for compounds that are on the sample summary report.

Regarding results on QC summary forms versus included raw data, due to calculations made at the instrument where many significant figures may be used, there may be slight discrepancies between the summary report result and that recorded on the raw data. This does not affect data usability.

VOLATILES

Regarding initial calibrations, it should be noted that the Quantitation Report concentrations supplied for the initial calibration reflect the calibration prior to updating. The response factors and areas are correct.

Regarding Quantitation Reports, it should be noted that the "#" symbol that appears on some of the Quantitation Reports is a software artifact and should be disregarded.

Compounds flagged with an "*" on the summary table have been calibrated using a non-average Response Factor calibration curve. The supporting curves are located after the initial calibration table.

Holding times were met for the samples.

Many of the samples had outliers for 1,2-Dichloroethane-d4 and 4-Bromofluorobenzene. All outliers have been flagged with an "*" on the summary form and the sample report.

Site specific QC was requested and analyzed on GEN-MW2-052322 and ethylbenzene recovered high in the matrix spike. This outlier has been flagged with an "*" on the summary page and an "M" on the sample report. Matrix interference is suspected. The Laboratory Control Sample recovered within acceptance limits.

The Method Blank was free from contamination within reportable ranges.

The instrument tunes passed all criteria and samples were within a 12-hour window.

The internal standards areas and retention times were within acceptance ranges for the samples and QC.

All data for the initial calibration was within acceptance limits for the reported analytes.

All continuing calibration data was within acceptance limits for the reported analytes.

Subcontracted Analyses

PAHs by EPA 8270D-SIM was sent to Alpha Analytical of Westborough, MA. Their reports are provided in their entirety as a separate entity after the Paradigm Environmental Services, Inc. report. Separate case narratives addressing the above parameters are included with their reports.

(signed) Steven DeVito Steven DeVito - Technical Director (date) 8/31/2022

BATCH LOG

Lab Name: <u>Paradigm Environmental Services</u>

Lab Project #: 222457
Client Name: Neu-Velle

Client Project Name: RGE Geneseo Fmr. MGP Site

Client Project #: N/A SDG No.: 2457-01

 Protocol:
 SW846
 Report Due Date:
 6/10/2022
 Batch Due Date:
 6/25/2022

LAB	MATRIX	CLIENT	REQUESTED ANALYSIS	DATE	DATE
SAMPLE NO.		SAMPLE ID	***	SAMPLED	REC'D
222457-01	Groundwater	GEN - MW3 - 052322	VOAs, SVOAs	5/23/2022	5/26/2022
222457-02	Groundwater	GEN - EB - 052322	VOAs, SVOAs	5/23/2022	5/26/2022
222457-03	Groundwater	GEN - DUP - 052322	VOAs, SVOAs	5/23/2022	5/26/2022
222457-04	Groundwater	GEN - MW2 - 052322	VOAs, SVOAs	5/23/2022	5/26/2022
222457-05	Groundwater	GEN - MW1 - 052422	VOAs, SVOAs	5/24/2022	5/26/2022
222457-06	Groundwater	GEN - MW4 - 052422	VOAs, SVOAs	5/24/2022	5/26/2022
222457-07	Groundwater	GEN - MW6 - 052522	VOAs, SVOAs	5/25/2022	5/26/2022
222457-08	Groundwater	GEN - MW8 - 052522	VOAs, SVOAs	5/25/2022	5/26/2022
222457-09	Groundwater	GEN - MW7 - 052622	VOAs, SVOAs	5/26/2022	5/26/2022
222457-10	Water	Trip Blank	VOAs	5/26/2022	5/26/2022

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CHAIN OF CUSTODY

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PARAI	DIGM

PARADIGM RGPROJECT REFERENCE FMN. MGP STT	CLIENT: NUI-VULL LADDRESS: 1667 LACK CITY COLLEGE STATE: N PHONE: 595 470 - 1 ATTN: Kgu N Matrix Codes: AQ - Aqueous Liquid NQ - Non-Aqueous Liquid	WG - Groundwater		Ema Kanasa SD - Soil SD - SL - Sludge PT -	Solid WP - Wipe	LUI — L 4 CON OL - Oil AR - Air
DATE COLLECTED TIME COLLECTED C O M P R C C C C O A F F C C C O B T T T E C C C O B T T T C C C C O B T T T C C C C C C C C C C C C C C C C	SAMPLE IDENTIFIER	CONTAINERS MAODES MAODES ACORES ACORES	12-to Sign		REMARKS	PARADIGM LAB SAMPLE NUMBER
5/23/22 14:00 × 5/23/22 12:45 × 5/23/22 × 5/23/22 × 5/24/22 15:30 × 5/24/22 16:30 × 5/25/22 14:45 × 5/25/22 16:30 × 5/25/22 16:30 × 5/26/22 13:00 × 5/20/2	GEN-MW3-05 GEN-EB-0523 GEN-DUP-052 GEN-MW2-05 GEN-MW1-05 GEN-MW6-05 GEN-MW8-05 GEN-MW3-05 GEN-MW3-05 GEN-MW3-05	2322 WG 3 X 22 WA 3 X 2322 WG 9 X 52422 WG 3 X 52422 WG 3 X 2522 WG 3 X 2522 WG 3 X 2522 WG 3 X 2522 WG 3 X	X	Equipo Pur MS		01 02 03 04 05 07 08
Turnaround Time Availability contingent upon lab app Standard 5 day None Required Batch QC Rush 3 day Category A Rush 2 day Rush 1 day	Report Supplements roval; additional fees may apply.	Received @ Lab By	Date/Time Date/Time Date/Time Date/Time	122 14:3	10°C ice S Total Cost: 1'4'6	
Oate Needed Other lease indicate date needed: please indicate paci	other EDD please indicate EDD needed :	Custody Seals By signing this form, clie	ent agrees to Paradigm	Terms and Condition	136 3 2 ns (reverse). page for sample cond	litions.

VOLATILE ORGANICS SAMPLE DATA



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW3 - 052322

Lab Sample ID:222457-01Date Sampled:5/23/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		6/2/2022 13:30
Ethylbenzene	< 2.00	ug/L		6/2/2022 13:30
m,p-Xylene	< 2.00	ug/L		6/2/2022 13:30
o-Xylene	< 2.00	ug/L		6/2/2022 13:30
Toluene	< 2.00	ug/L		6/2/2022 13:30

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4	83.7	81.1 - 136		6/2/2022	13:30
4-Bromofluorobenzene	87.0	75.8 - 132		6/2/2022	13:30
Pentafluorobenzene	121	82 - 132		6/2/2022	13:30
Toluene-D8	115	64.6 - 137		6/2/2022	13:30

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09594.D



Client: **Neu-Velle**

RGE Geneseo Fmr. MGP Site **Project Reference:**

Sample Identifier: GEN - EB - 052322

Date Sampled: Lab Sample ID: 222457-02 5/23/2022 **Date Received:** 5/26/2022 **Matrix:** Groundwater

Volatile Organics

Analyte	Result	<u>Units</u>		Qualifier	Date A	nalyzed
Benzene	< 1.00	ug/L		UJ	6/2/20	22 13:49
Ethylbenzene	< 2.00	ug/L			6/2/20	22 13:49
m,p-Xylene	< 2.00	ug/L			6/2/20	22 13:49
o-Xylene	< 2.00	ug/L			6/2/20	22 13:49
Toluene	< 2.00	ug/L		\downarrow	6/2/20	22 13:49
<u>Surrogate</u>	<u>Percen</u>	t Recovery	<u>Limits</u>	<u>Outliers</u>	Date An	alyzed
1,2-Dichloroethane-d4	7	74.1	81.1 - 136	*	6/2/2022	13:49
4-Bromofluorobenzene	7	76.6	75.8 - 132		6/2/2022	13:49
Pentafluorobenzene		126	82 - 132		6/2/2022	13:49
Toluene-D8	:	106	64.6 - 137		6/2/2022	13:49

Method Reference(s): EPA 8260C EPA 5030C

Data File: z09595.D

> MKP 9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - DUP - 052322

Lab Sample ID:222457-03Date Sampled:5/23/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>		Qualifier	Date A	nalyzed
Benzene	< 1.00	ug/L		UJ	6/2/20	22 14:09
Ethylbenzene	< 2.00	ug/L			6/2/20	22 14:09
m,p-Xylene	< 2.00	ug/L			6/2/20	22 14:09
o-Xylene	< 2.00	ug/L			6/2/20	22 14:09
Toluene	< 2.00	ug/L		lacksquare	6/2/20	22 14:09
<u>Surrogate</u>	<u>Per</u>	cent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		72.5	81.1 - 136	*	6/2/2022	14:09
4-Bromofluorobenzene		74.3	75.8 - 132	*	6/2/2022	14:09
Pentafluorobenzene		129	82 - 132		6/2/2022	14:09
Toluene-D8		109	64.6 - 137		6/2/2022	14:09
Method Reference(s):	EPA 8260C					

Method Reference(s): EPA 8260C EPA 5030C

Data File: z09596.D

MKP 9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW2 - 052322

Lab Sample ID:222457-04Date Sampled:5/23/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date A	nalyzed
Benzene	< 1.00	ug/L	U	J	6/2/20	22 14:28
Ethylbenzene	< 2.00	ug/L		M	6/2/20	22 14:28
m,p-Xylene	< 2.00	ug/L			6/2/20	22 14:28
o-Xylene	< 2.00	ug/L			6/2/20	22 14:28
Toluene	< 2.00	ug/L			6/2/20	22 14:28
<u>Surrogate</u>	<u>Perce</u>	nt Recovery	<u>Limits</u>	Outliers	Date An	<u>alyzed</u>
1,2-Dichloroethane-d4		70.5	81.1 - 136	*	6/2/2022	14:28
4-Bromofluorobenzene		68.1	75.8 - 132	*	6/2/2022	14:28
Pentafluorobenzene		126	82 - 132		6/2/2022	14:28
Toluene-D8		100	64.6 - 137		6/2/2022	14:28

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09597.D

MKP 9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW1 - 052422

Lab Sample ID:222457-05Date Sampled:5/24/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date A	<u>nalyzed</u>
Benzene	< 1.00	ug/L		UJ	6/2/20	22 15:45
Ethylbenzene	< 2.00	ug/L			6/2/20	22 15:45
m,p-Xylene	< 2.00	ug/L			6/2/20	22 15:45
o-Xylene	< 2.00	ug/L			6/2/20	22 15:45
Toluene	< 2.00	ug/L		\downarrow	6/2/20	22 15:45
<u>Surrogate</u>	<u>Perc</u>	ent Recovery	<u>Limits</u>	Outliers	<u>Date An</u>	<u>ialyzed</u>
1,2-Dichloroethane-d4		72.8	81.1 - 136	*	6/2/2022	15:45
4-Bromofluorobenzene		72.2	75.8 - 132	*	6/2/2022	15:45
Pentafluorobenzene		125	82 - 132		6/2/2022	15:45
Toluene-D8		106	64.6 - 137		6/2/2022	15:45

Method Reference(s): EPA 8260C EPA 5030C Data File: z09601.D

MKP

9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW4 - 052422

Lab Sample ID:222457-06Date Sampled:5/24/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

Analyte	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	0.503	ug/L	J J	6/2/2022 16:05
Ethylbenzene	< 2.00	ug/L	UJ	6/2/2022 16:05
m,p-Xylene	2.98	ug/L	J	6/2/2022 16:05
o-Xylene	< 2.00	ug/L	UJ	6/2/2022 16:05
Toluene	< 2.00	ug/L	UJ	6/2/2022 16:05

<u>Surrogate</u>	Percent Recovery	<u>Limits</u>	<u>Outliers</u>	<u>Date Ar</u>	<u>ıalyzed</u>
1,2-Dichloroethane-d4	68.6	81.1 - 136	*	6/2/2022	16:05
4-Bromofluorobenzene	68.9	75.8 - 132	*	6/2/2022	16:05
Pentafluorobenzene	122	82 - 132		6/2/2022	16:05
Toluene-D8	98.6	64.6 - 137		6/2/2022	16:05

Method Reference(s): EPA 8260C

EPA 5030C **Data File:** z09602.D

MKP 9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW6 - 052522

Lab Sample ID:222457-07Date Sampled:5/25/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date A	<u>nalyzed</u>
Benzene	164	ug/L		J	6/2/20	22 15:26
Ethylbenzene	46.7	ug/L			6/2/20	22 15:26
m,p-Xylene	55.7	ug/L			6/2/20	22 15:26
o-Xylene	48.5	ug/L			6/2/20	22 15:26
Toluene	49.8	ug/L		\downarrow	6/2/20	22 15:26
<u>Surrogate</u>	<u>Per</u>	cent Recovery	<u>Limits</u>	Outliers	Date An	alyzed
1,2-Dichloroethane-d4		79.8	81.1 - 136	*	6/2/2022	15:26
4-Bromofluorobenzene		88.8	75.8 - 132		6/2/2022	15:26
Pentafluorobenzene		122	82 - 132		6/2/2022	15:26
Toluene-D8		108	64.6 - 137		6/2/2022	15:26

Method Reference(s): EPA 8260C EPA 5030C

Data File: z09600.D

MKP 9/8/2022



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW8 - 052522

Lab Sample ID:222457-08Date Sampled:5/25/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	<u>Result</u>	<u>Units</u>		Qualifier	Date A	<u>ıalyzed</u>
Benzene	1.88	ug/L	J		6/2/20	22 16:24
Ethylbenzene	3.08	ug/L			6/2/20	22 16:24
m,p-Xylene	5.86	ug/L			6/2/20	22 16:24
o-Xylene	3.12	ug/L			6/2/20	22 16:24
Toluene	2.00	ug/L	$lack \psi$	J	6/2/20	22 16:24
Surrogate	<u>Perc</u>	ent Recovery	<u>Limits</u>	<u>Outliers</u>	Date An	<u>alyzed</u>
1,2-Dichloroethane-d4		75.4	81.1 - 136	*	6/2/2022	16:24
4-Bromofluorobenzene		79.3	75.8 - 132		6/2/2022	16:24
Pentafluorobenzene		121	82 - 132		6/2/2022	16:24
Toluene-D8		106	64.6 - 137		6/2/2022	16:24

Method Reference(s): EPA 8260C EPA 5030C Data File: z09603.D

9/8/2022

MKP



Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: GEN - MW7 - 052622

Lab Sample ID:222457-09Date Sampled:5/26/2022Matrix:GroundwaterDate Received:5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>	Qualifier	Date Analyzed
Benzene	< 1.00	ug/L		6/2/2022 16:43
Ethylbenzene	< 2.00	ug/L		6/2/2022 16:43
m,p-Xylene	< 2.00	ug/L		6/2/2022 16:43
o-Xylene	< 2.00	ug/L		6/2/2022 16:43
Toluene	< 2.00	ug/L		6/2/2022 16:43

Surrogate	Percent Recovery	<u>Limits</u>	Outliers	Date Analyzed	
1,2-Dichloroethane-d4	81.6	81.1 - 136		6/2/2022	16:43
4-Bromofluorobenzene	78.8	75.8 - 132		6/2/2022	16:43
Pentafluorobenzene	113	82 - 132		6/2/2022	16:43
Toluene-D8	103	64.6 - 137		6/2/2022	16:43

Method Reference(s): EPA 8260C

EPA 5030C

Data File: z09604.D



5/26/2022

Date Sampled:

Client: <u>Neu-Velle</u>

Project Reference: RGE Geneseo Fmr. MGP Site

Sample Identifier: Trip Blank
Lab Sample ID: 222457-10

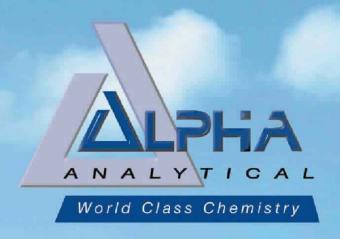
Matrix: Water Date Received: 5/26/2022

Volatile Organics

<u>Analyte</u>	Result	<u>Units</u>		Qualifier	Date Ar	<u>ıalyzed</u>
Benzene	< 1.00	ug/L		UJ	6/2/20	22 12:07
Ethylbenzene	< 2.00	ug/L			6/2/20	22 12:07
m,p-Xylene	< 2.00	ug/L			6/2/20	22 12:07
o-Xylene	< 2.00	ug/L			6/2/20	22 12:07
Toluene	< 2.00	ug/L		\downarrow	6/2/20	22 12:07
<u>Surrogate</u>	Percent Recovery		<u>Limits</u>	<u>Outliers</u>	Date Analyzed	
1,2-Dichloroethane-d4		75.8	81.1 - 136	*	6/2/2022	12:07
4-Bromofluorobenzene		78.7	75.8 - 132		6/2/2022	12:07
Pentafluorobenzene		121	82 - 132		6/2/2022	12:07
Toluene-D8		107	64.6 - 137		6/2/2022	12:07

Method Reference(s): EPA 8260C
EPA 5030C
Data File: z09590.D

MKP 9/8/2022



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Alpha Analytical

Laboratory Code: 11148

SDG Number: L2228058

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 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2228058-01	GEN MW3-052322 222457-01	WATER	Not Specified	05/23/22 14:00	05/26/22
L2228058-02	GEN-EB-052322 222457-02	WATER	Not Specified	05/23/22 12:45	05/26/22
L2228058-03	GEN-DUP-052322 222457-03	WATER	Not Specified	05/23/22 00:00	05/26/22
L2228058-04	GEN-MW2-052322 222457-04	WATER	Not Specified	05/23/22 15:30	05/26/22
L2228058-05	GEN-MW1-052422 222457-05	WATER	Not Specified	05/24/22 14:45	05/26/22
L2228058-06	GEN-MW4-052422 222457-06	WATER	Not Specified	05/24/22 16:30	05/26/22
L2228058-07	GEN-MW6-052522 222457-07	WATER	Not Specified	05/25/22 14:45	05/26/22
L2228058-08	GEN-MW8-052522 222457-08	WATER	Not Specified	05/25/22 16:30	05/26/22
L2228058-09	GEN-MW7-052622 222457-09	WATER	Not Specified	05/26/22 13:00	05/26/22



 Project Name:
 222457
 Lab Number:
 L2228058

 Project Number:
 222457
 Report Date:
 06/09/22

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature: Wallin Wallier Report Date: 06/09/22

Title: Technical Director/Representative





PARADIGM

CHAIN OF CUSTODY L2228058

FA	RADIGM	COMPA	Paradigm Environme	ental	COMPAN	IÝ:	Same		LAB PROJECT #: C	LIENT PROJECT #:	
1	ss: 179 Lake Avenue	1000	ADDRES	S:							
1		CITY:	Rochester STATE: N	Y ZIP: 14608	CITY		STATE:	ZIP:	TURNAROUND TIME: (WORK	(ING DAYS)	
100	-	PHONE	FAX:	147/	PHONE:		FAX:			STD	OTHER
PROJECT NAME/SI	TE NAME:	ATTN:	Reporting		ATTN:	Ac	counts Payable				0 Day
		COMME	NTS: Please email results to	o reporting@pa	aradigi	menv.c	om				550,
	a systematic	OF THE PERSON	CONTRACTOR OF THE PARTY			RE	QUESTED ANALYSIS	HT SING	Date Due: 6/10/	22 for data	Manual Indian
DATE	TIME OS		SAMPLE LOCATION/FIELD ID	M A T R I	C O N T A I N E R E R	8270 PAH		4	REMARKS	PARADIGM LAB SAME	PLE NUMBER
1 5/23/22	14:00	X	GEN-MW3-052322	WG	1	X	NYS ASP Cat	NYS ASP Cat B Report Pkg and			
2 5/23/22	12:45	X	GEN-EB-052322	WG	1	X	Report Pkg an				
3 5/23/22		X	GEN-DUP-052322	WG	1	X	NYSDEC EDD		03		
5/23/22	15:30	X	GEN-MW2-052322	WG	1	x	SIM for PAH DLs	SIM for PAH DLs as needed			
5/24/22	14:45	X	GEN-MW1-052422	WG	1	x	needed				
5/24/22	16:30	X	GEN-MW4-052422	WG	1	х					
5/25/22	14:45	Х	GEN-MW6-052522	WG	্ৰ	x	Run MS/MSD o	,	07		
5/25/22	16:30	×	GEN-MW8-052522	WG	1	x	sample GEN-MV		08		
5/26/22	13:00	X	GEN-MW7-052622	WG	1	x	052322		09		
0											
	ONLY BELOW TH on: Per NELAC/ELAP Receipt Parameter Container Type: Preservation:		Y N N	Relinquished B	lai	1	5/2 6/2-2 Date/Time	1600	Total Cos	t	
omments.	Holding Time: Temperature:		Y N	Received By	ny	Kon	1 19.4 L 5/24 Date/Time	122 0	/6:00 P.I.F.		
				Received @ Lat	Ву		Date/Time				

Semivolatiles Data by Method 8270D-SIM

Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-01

Client ID : GEN MW3-052322 222457-01

Sample Location :

GPC Cleanup

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-01
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

: N

Date Received : 05/26/22
Date Analyzed : 05/29/22 16:19
Date Extracted : 05/27/22
Dilution Factor : 1

Project Number : 222457

Date Collected : 05/23/22 14:00

: L2228058

Lab Number

Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	UJ
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	
91-57-6	2-Methylnaphthalene	ND	0.10	0.05	U	



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-02

Client ID : GEN-EB-052322 222457-02

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-02
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058

Project Number : 222457

Date Collected : 05/23/22 12:45

Date Received : 05/26/22 Date Analyzed : 05/29/22 16:35

Date Extracted : 05/27/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L			
CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	UJ
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	
91-57-6	2-Methylnaphthalene	ND	0.10	0.05	U	
	7 - F					



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-03

Client ID : GEN-DUP-052322 222457-03

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-03
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058 Project Number : 222457

Date Collected : 05/23/22 00:00
Date Received : 05/26/22

Date Analyzed : 05/29/22 16:52

Date Extracted : 05/27/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
83-32-9	Acenaphthene	ND	0.10	0.04	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U
206-44-0	Fluoranthene	ND	0.10	0.04	U
91-20-3	Naphthalene	ND	0.10	0.04	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U UJ
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U
218-01-9	Chrysene	ND	0.10	0.04	U
208-96-8	Acenaphthylene	ND	0.10	0.04	U
120-12-7	Anthracene	ND	0.10	0.04	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U
86-73-7	Fluorene	ND	0.10	0.04	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U
129-00-0	Pyrene	ND	0.10	0.04	U
1-57-6	2-Methylnaphthalene	ND	0.10	0.05	U
-					



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-04

Client ID : GEN-MW2-052322 222457-04

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-04
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058
Project Number : 222457
Posto Colleget d : 05/23/23 18

Date Collected : 05/23/22 15:30 Date Received : 05/26/22

Date Analyzed : 05/29/22 17:08 Date Extracted : 05/27/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L			
ım	eter	Results	RL	MDL	Qualifie	r
ap	phthene	ND	0.10	0.04	U	
lo	ronaphthalene	ND	0.20	0.04	U	
raı	nthene	ND	0.10	0.04	U	
nth	alene	ND	0.10	0.04	U	
20(a)anthracene	ND	0.10	0.02	U	
20(a)pyrene	ND	0.10	0.04	U	UJ
20(b)fluoranthene	ND	0.10	0.02	U	
20(k)fluoranthene	ND	0.10	0.04	U	UJ
se	ne	ND	0.10	0.04	U	
ıaı	ohthylene	ND	0.10	0.04	U	
ra	cene	ND	0.10	0.04	U	
20(ghi)perylene	ND	0.10	0.04	U	UJ
rei	ne	ND	0.10	0.04	U	
naı	nthrene	ND	0.10	0.02	U	
nz	o(a,h)anthracene	ND	0.10	0.04	U	UJ
no	(1,2,3-cd)pyrene	ND	0.10	0.04	U	UJ
ne		ND	0.10	0.04	U	
th	ylnaphthalene	ND	0.10	0.05	U	
ne		ND	0.10	0.04		U



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-05

Client ID : GEN-MW1-052422 222457-05

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-05
Sample Amount : 1000 ml
Extraction Method : EPA 3510C

Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058

Project Number : 222457

Date Collected : 05/24/22 14:45

Date Received : 05/26/22

Date Analyzed : 05/29/22 18:46

Date Extracted : 05/27/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L		
CAS NO.	Parameter	Results	RL	MDL	Qualifier
83-32-9	Acenaphthene	ND	0.10	0.04	U
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U
206-44-0	Fluoranthene	ND	0.10	0.04	U
91-20-3	Naphthalene	ND	0.10	0.04	U
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	u UJ
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U
218-01-9	Chrysene	ND	0.10	0.04	U
208-96-8	Acenaphthylene	ND	0.10	0.04	U
120-12-7	Anthracene	ND	0.10	0.04	U
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U
86-73-7	Fluorene	ND	0.10	0.04	U
85-01-8	Phenanthrene	ND	0.10	0.02	U
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U
129-00-0	Pyrene	ND	0.10	0.04	U
91-57-6	2-Methylnaphthalene	ND	0.10	0.05	U



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-06

Client ID : GEN-MW4-052422 222457-06

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-06
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058

Project Number : 222457

Date Collected : 05/24/22 16:30

Date Received : 05/26/22

Date Analyzed : 05/29/22 19:02

Date Extracted : 05/27/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

			ug/L				
CAS NO.	Parameter	Results	RL	MDL	Qualifier		
83-32-9	Acenaphthene	ND	0.10	0.04	U		
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U		
206-44-0	Fluoranthene	ND	0.10	0.04	U		
91-20-3	Naphthalene	ND	0.10	0.04	U		
56-55-3	Benzo(a)anthracene	ND	0.10	0.02	U		
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U UJ		
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U		
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U		
218-01-9	Chrysene	ND	0.10	0.04	U		
208-96-8	Acenaphthylene	ND	0.10	0.04	U		
120-12-7	Anthracene	ND	0.10	0.04	U		
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U		
86-73-7	Fluorene	ND	0.10	0.04	U		
85-01-8	Phenanthrene	0.03	0.10	0.02	J		
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U		
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U		
129-00-0	Pyrene	ND	0.10	0.04	U		
91-57-6	2-Methylnaphthalene	ND	0.10	0.05	U		



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-07

Client ID : GEN-MW6-052522 222457-07

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-07
Sample Amount : 1000 ml
Extraction Method : EPA 3510C

Extract Volume : 1000 uL GPC Cleanup : N

Lab Number : L2228058
Project Number : 222457
Date Collected : 05/25/22 14:45

Date Received : 05/26/22 Date Analyzed : 05/31/22 20:01

Date Extracted : 05/29/22

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

				ug/L			
CAS NO.	Parameter		Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene		2.9	0.10	0.04		
91-58-7	2-Chloronaphthalene		ND	0.20	0.04	U	
206-44-0	Fluoranthene		0.05	0.10	0.04	J	
91-20-3	Naphthalene		200	0.10	0.04	E	
56-55-3	Benzo(a)anthracene	0.05 UJ	0.02	0.10	0.02	J	
50-32-8	Benzo(a)pyrene		ND	0.10	0.04	U	UJ
205-99-2	Benzo(b)fluoranthene		ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene		ND	0.10	0.04	U	
218-01-9	Chrysene		ND	0.10	0.04	U	
208-96-8	Acenaphthylene		29	0.10	0.04	E	
120-12-7	Anthracene		0.42	0.10	0.04		
191-24-2	Benzo(ghi)perylene		ND	0.10	0.04	U	
86-73-7	Fluorene		5.5	0.10	0.04		
85-01-8	Phenanthrene		1.9	0.10	0.02		
53-70-3	Dibenzo(a,h)anthracene		ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene		ND	0.10	0.04	U	UJ
129-00-0	Pyrene		ND	0.10	0.04	U	
91-57-6	2-Methylnaphthalene		0.10	0.10	0.05	J	
	<u> </u>						



Client : Paradigm Environmental Services Lab Number : L2228058
Project Name : 222457 Project Number : 222457

Sample Matrix Date Extracted : WATER : 05/29/22 Analytical Method : 1,8270D-SIM **Dilution Factor** : 10 Lab File ID : 58-07D1 Analyst : 1000 ml Sample Amount Instrument ID : SV118 Extraction Method : EPA 3510C GC Column : RXI-5SilM **Extract Volume** : 1000 uL %Solids : N/A

GPC Cleanup : N Injection Volume : 1 uL

ug/L CAS NO. Results RL MDL Qualifier **Parameter** 91-20-3 180 Naphthalene 1.0 0.43 21 208-96-8 Acenaphthylene 1.0 0.35



Client : Paradigm Environmental Services

Project Name : 222457

Lab ID : L2228058-08

Client ID : GEN-MW8-052522 222457-08

Sample Location :

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-08
Sample Amount : 1000 ml
Extraction Method : EPA 3510C
Extract Volume : 1000 uL

GPC Cleanup : N

Lab Number : L2228058
Project Number : 222457
Page Collected : 05/05/32.1

Date Collected : 05/25/22 16:30 Date Received : 05/26/22

Date Analyzed : 05/31/22 20:17

Date Extracted : 05/29/22
Dilution Factor : 1
Analyst : RP

Injection Volume : 1 uL

Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

ug/L Results MDL Qualifier CAS NO. **Parameter** RL 83-32-9 Acenaphthene ND 0.10 0.04 U 91-58-7 2-Chloronaphthalene ND 0.20 0.04 U 206-44-0 Fluoranthene ND 0.10 0.04 U 91-20-3 Naphthalene ND 0.10 0.04 U 56-55-3 Benzo(a)anthracene 0.10 UJ 0.02 0.10 0.02 J 50-32-8 Benzo(a)pyrene ND 0.10 0.04 U UJ U 205-99-2 Benzo(b)fluoranthene ND 0.10 0.02 207-08-9 Benzo(k)fluoranthene ND 0.10 0.04 U U 218-01-9 Chrysene ND 0.10 0.04 0.04 U 208-96-8 Acenaphthylene ND 0.10 0.10 0.04 U 120-12-7 Anthracene ND 191-24-2 Benzo(ghi)perylene ND 0.10 0.04 п 86-73-7 Fluorene ND 0.10 0.04 U 85-01-8 Phenanthrene ND 0.10 0.02 U 53-70-3 Dibenzo(a,h)anthracene ND 0.10 0.04 U UJ 193-39-5 ND 0.10 0.04 U Indeno(1,2,3-cd)pyrene 129-00-0 Pyrene ND 0.10 0.04 U 91-57-6 ND 0.10 0.05 2-Methylnaphthalene U



Client : Paradigm Environmental Services

Project Name : 222457 Lab ID : L2228058-09

Client ID : GEN-MW7-052622 222457-09

Sample Location :

GPC Cleanup

Sample Matrix : WATER
Analytical Method : 1,8270D-SIM
Lab File ID : 28058-09
Sample Amount : 900 ml
Extraction Method : EPA 3510C
Extract Volume : 900 uL

: N

Project Number : 222457

Date Collected : 05/26/22 13:00

Date Received : 05/26/22

Date Analyzed : 05/31/22 20:34

Date Extracted : 05/29/22

: L2228058

Dilution Factor : 1
Analyst : RP
Instrument ID : SV115
GC Column : RXI-5SilM
%Solids : N/A

Injection Volume : 1 uL

Lab Number

Parameter Acenaphthene 2-Chloronaphthalene	Results ND ND	0.11	MDL 0.04	Qualifier U
2-Chloronaphthalene		0.11	0.04	
2-Chloronaphthalene		0.11	0.04	11
·	ND			<u> </u>
= 1 .1	ND	0.22	0.04	U
Fluoranthene	ND	0.11	0.04	U
Naphthalene	ND	0.11	0.05	U
Benzo(a)anthracene	ND	0.11	0.02	U
Benzo(a)pyrene	ND	0.11	0.04	U UJ
Benzo(b)fluoranthene	ND	0.11	0.02	U
Benzo(k)fluoranthene	ND	0.11	0.05	U
Chrysene	ND	0.11	0.04	U
Acenaphthylene	ND	0.11	0.04	U
Anthracene	ND	0.11	0.04	U
Benzo(ghi)perylene	ND	0.11	0.05	U
Fluorene	ND	0.11	0.04	U
Phenanthrene	0.03	0.11	0.02	J
Dibenzo(a,h)anthracene	ND	0.11	0.04	U
Indeno(1,2,3-cd)pyrene	ND	0.11	0.04	U UJ
Pyrene	ND	0.11	0.04	U
2-Methylnaphthalene	ND	0.11	0.05	U
	Benzo(a)anthracene Benzo(b)fluoranthene Benzo(k)fluoranthene Chrysene Acenaphthylene Anthracene Benzo(ghi)perylene Fluorene Phenanthrene Dibenzo(a,h)anthracene Indeno(1,2,3-cd)pyrene Pyrene	Naphthalene ND Benzo(a)anthracene ND Benzo(a)pyrene ND Benzo(b)fluoranthene ND Benzo(k)fluoranthene ND Chrysene ND Acenaphthylene ND Anthracene ND Benzo(ghi)perylene ND Fluorene ND Phenanthrene 0.03 Dibenzo(a,h)anthracene ND Indeno(1,2,3-cd)pyrene ND Pyrene ND	Naphthalene ND 0.11 Benzo(a)anthracene ND 0.11 Benzo(a)pyrene ND 0.11 Benzo(b)fluoranthene ND 0.11 Benzo(k)fluoranthene ND 0.11 Chrysene ND 0.11 Acenaphthylene ND 0.11 Anthracene ND 0.11 Benzo(ghi)perylene ND 0.11 Fluorene ND 0.11 Phenanthrene 0.03 0.11 Dibenzo(a,h)anthracene ND 0.11 Indeno(1,2,3-cd)pyrene ND 0.11 Pyrene ND 0.11	Naphthalene ND 0.11 0.05 Benzo(a)anthracene ND 0.11 0.02 Benzo(a)pyrene ND 0.11 0.04 Benzo(b)fluoranthene ND 0.11 0.02 Benzo(k)fluoranthene ND 0.11 0.05 Chrysene ND 0.11 0.04 Acenaphthylene ND 0.11 0.04 Anthracene ND 0.11 0.04 Benzo(ghi)perylene ND 0.11 0.05 Fluorene ND 0.11 0.04 Phenanthrene 0.03 0.11 0.02 Dibenzo(a,h)anthracene ND 0.11 0.04 Indeno(1,2,3-cd)pyrene ND 0.11 0.04 Pyrene ND 0.11 0.04



Appendix B

Laboratory QC Documentation

2 VOLATILE SURROGATE RECOVERY

Lab Name: <u>Paradigm Environmental Services</u>

Lab Project #: 222457
Client Name: Neu-Velle

Client Project Name: RGE Geneseo Fmr. MGP Site

Client Project #: <u>N/A</u> SDG No.: <u>2457-01</u>

Matrix: Groundwater
QC Batch: voaw220602

Instrument ID: <u>Instrument1</u>

GC Column 1: DB-624 ID (mm): 0.20 Detector: MSD

	LAB	CLIENT	PFB	12DCEd4	TD8	4BFB	Total
	SAMPLE NO.	SAMPLE ID	%REC	%REC	%REC	%REC	Out
	Blk 1	N/A	116	82.6	107	79.3	0
	LCS 1	N/A	116	67.0 *	97.8	75.3 *	2
3	222457-01	GEN - MW3 - 052322	121	83.7	115	87.0	0
4	222457-02	GEN - EB - 052322	126	74.1 *	106	76.6	1
5	222457-03	GEN - DUP - 052322	129	72.5 *	109	74.3 *	2
6	222457-04	GEN - MW2 - 052322	126	70.5 *	100	68.1 *	2
7	222457-04MS	GEN - MW2 - 052322 MS	123	75.1 *	108	89.5	1
8	222457-04MSD	GEN - MW2 - 052322 MSD	122	86.6	110	95.7	0
9	222457-05	GEN - MW1 - 052422	125	72.8 *	106	72.2 *	2
10	222457-06	GEN - MW4 - 052422	122	68.6 *	98.6	68.9 *	2
11	222457-07	GEN - MW6 - 052522	122	79.8 *	108	88.8	1
12	222457-08	GEN - MW8 - 052522	121	75.4 *	106	79.3	1
13	222457-09	GEN - MW7 - 052622	113	81.6	103	78.8	0
14	222457-10	Trip Blank	121	75.8 *	107	78.7	1
15							
16							
17							
18							
19							
20							
21							
22							
23							
24							
25							

QC LIMITS %
(82 - 132)
(81.1 - 136)
(64.6 - 137)
(75.8 - 132)

^{*} Values outside of current required QC limits

D Surrogate diluted out



QC Report for Matrix Spike and Matrix Spike Duplicate

SDG #: 2457-01

Client: Neu-Velle Lab Project ID: 222457

Project Reference: RGE Geneseo Fmr. MGP Site

Lab Sample ID: 222457-04

Sample Identifier: GEN - MW2 - 052322

Matrix: Groundwater

Date Sampled: 5/23/2022 **Date Received:** 5/26/2022

Date Analyzed: 6/2/2022

Volatile Organics

	Sample	Result	<u>MS</u>	<u>MS</u>	MS %	MSD	MSD	MSD %	<u>% Rec.</u>	<u>MS</u>	<u>MSD</u>	<u>Relative</u>	<u>RPD</u>	<u>RPD</u>
<u>Analyte</u>	Result	<u>Units</u>	Added	Result	Recovery	Added	Result	Recovery	<u>Limits</u>	<u>Outlier</u>	Outlier	% Diff.	Limit	<u>Outlier</u>
Benzene	< 1.00	ug/L	50.0	56.4	113	50.0	53.7	107	81.6 - 114			4.81	15	
Ethylbenzene	< 2.00	ug/L	50.0	57.7	115	50.0	52.6	105	72.1 - 110	*		9.27	20	
Toluene	< 2.00	ug/L	50.0	58.2	116	50.0	56.4	113	62.9 - 125			3.11	32	

Method Reference(s): EPA 8260C

EPA 5030C

Data File(s): z09598.D

z09599.D z09597.D

1

QC Batch ID: voaw220602

Any estimated values are displayed, and derived values calculated, based on numeric result only. See primary analytical report for data flags.

This report is part of a multipage document and should only be evaluated in its entirety. The Chain of Custody provides additional sample information, including compliance with the sample condition requirements upon receipt.

Matrix Spike Sample Summary Form 3 **Semivolatiles**

Client : Paradigm Environmental Services

Project Name : 222457

Client Sample ID : GEN-MW2-052322 222457-04

Lab Sample ID : L2228058-04 Matrix Spike : WG1643559-4 Matrix Spike Dup : WG1643559-5

Lab Number : L2228058 Project Number : 222457

Matrix : WATER
Analysis Date : 05/29/22 17:08 MS Analysis Date : 06/09/22 17:48 MSD Analysis Date: 06/09/22 18:04

		Matrix Spike Sample			Matrix Spike Duplicate					
	Sample	Spike	Spike		Spike	Spike				
	Conc.	Added	Conc.	%R	Added	Conc.	%R	RPD	Recovery	RPD
Parameter	(ug/l)	(ug/l)	(ug/l)		(ug/l)	(ug/l)			Limits	Limit
Acenaphthene	ND	10	6.4	64	10	7.8	78	20	40-140	40
2-Chloronaphthalene	ND	10	6.3	63	10	7.7	77	20	40-140	40
Fluoranthene	ND	10	7.8	78	10	9.3	93	18	40-140	40
Naphthalene	ND	10	6.2	62	10	7.5	75	19	40-140	40
Benzo(a)anthracene	ND	10	7.0	70	10	8.5	85	19	40-140	40
Benzo(a)pyrene	ND	10	4.4	44	10	5.0	50	13	40-140	40
Benzo(b)fluoranthene	ND	10	4.2	42	10	5.0	50	17	40-140	40
Benzo(k)fluoranthene	ND	10	3.8	38	10	4.2	42	10	40-140	40
Chrysene	ND	10	5.3	53	10	6.2	62	16	40-140	40
Acenaphthylene	ND	10	6.9	69	10	8.4	84	20	40-140	40
Anthracene	ND	10	6.7	67	10	8.1	81	19	40-140	40
Benzo(ghi)perylene	ND	10	1.2	12	10	1.4	14 Q	15	40-140	40
Fluorene	ND	10	6.8	68	10	8.2	82	19	40-140	40
Phenanthrene	ND	10	6.1	61	10	7.5	75	21	40-140	40
Dibenzo(a,h)anthracene	ND	10	1.4	14 Q	10	1.6	16 Q	13	40-140	40
Indeno(1,2,3-cd)pyrene	ND	10	1.4	14 0	10	1.7	17 Q	19	40-140	40
Pyrene	ND	10	8.1	81	10	9.5	95	16	40-140	40
2-Methylnaphthalene	ND	10	6.2	62	10	7.5	75	19	40-140	40



Client : Paradigm Environmental Services

Project Name : 222457

Lab ID : WG1644257-1

Client ID : WG1644257-1BLANK

Sample Location :

Sample Matrix : WATER Analytical Method : 1,8270D-SIM Lab File ID : 644257-1 Sample Amount : 1000 ml Extraction Method : EPA 3510C

Extract Volume : 1000 uL GPC Cleanup : N

Lab Number : L2228058 Project Number : 222457 Date Collected : NA Date Received : NA Date Analyzed

: 05/31/22 19:45

Date Extracted : 05/29/22

Dilution Factor : 1 Analyst : RP Instrument ID : SV115 GC Column : RXI-5SilM %Solids : N/A

Injection Volume : 1 uL

CAS NO.	Parameter	Results	RL	MDL	Qualifier	
83-32-9	Acenaphthene	ND	0.10	0.04	U	
91-58-7	2-Chloronaphthalene	ND	0.20	0.04	U	
206-44-0	Fluoranthene	ND	0.10	0.04	U	
91-20-3	Naphthalene	ND	0.10	0.04	U	
56-55-3	Benzo(a)anthracene	0.02	0.10	0.02	J	
50-32-8	Benzo(a)pyrene	ND	0.10	0.04	U	
205-99-2	Benzo(b)fluoranthene	ND	0.10	0.02	U	
207-08-9	Benzo(k)fluoranthene	ND	0.10	0.04	U	
218-01-9	Chrysene	ND	0.10	0.04	U	
208-96-8	Acenaphthylene	ND	0.10	0.04	U	
120-12-7	Anthracene	ND	0.10	0.04	U	
191-24-2	Benzo(ghi)perylene	ND	0.10	0.04	U	
86-73-7	Fluorene	ND	0.10	0.04	U	
85-01-8	Phenanthrene	ND	0.10	0.02	U	
53-70-3	Dibenzo(a,h)anthracene	ND	0.10	0.04	U	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.10	0.04	U	
129-00-0	Pyrene	ND	0.10	0.04	U	
91-57-6	2-Methylnaphthalene	ND	0.10	0.05	U	



Calibration Verification Summary Form 7 Semivolatiles

Client : Paradigm Environmental Services Lab Number : L2228058
Project Name : 222457 Project Number : 222457

Project Name : 222457 Project Number : 222457 Calibration Date : 05/29/22 07:24

 Lab File ID
 : CCV0529A
 Init. Calib. Date(s)
 : 02/04/22

 Sample No
 : WG1644230-3
 Init. Calib. Times
 : 15:41
 18:08

Channel :

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
1,4-Dichlorobenzene-d4	1	1	.05	0	20	100	0
2-Fluorophenol	0.944	1.021	.05	-8.2	20	104	0
Phenol-d6	1.196	1.289	.05	-7.8	20	101	0
Hexachloroethane	0.496	0.531	.05	-7.1	20	103	0
Bis(2-chloroethyl)ether	1.127	1.097	.05	2.7	20	95	0
n-nitrosodi-n-propylamine	0.764	0.857	.05	-12.2	20	104	0
Nitrobenzene-d5	0.917	1.295	.05	-41.2*	20	127	0
Naphthalene-d8	1	1	.05	0	20	105	0
Naphthalene	1.025	0.953	.05	7	20	98	0
Hexachlorobutadiene	0.189	0.182	.05	3.7	20	102	0
2-Methylnaphthalene	0.627	0.611	.05	2.6	20	101	0
1-Methylnaphthalene	0.607	0.597	.05	1.6	20	105	0
2-Fluorobiphenyl	0.735	0.688	.05	6.4	20	106	0
2-Chloronaphthalene	0.649	0.632	.05	2.6	20	103	0
Acenaphthylene	0.972	1.102	.05	-13.4	20	112	0
Acenaphthene-d10	1	1	.05	0	20	118	0
Acenaphthene	1.325	1.202	.05	9.3	20	107	0
Fluorene	1.404	1.392	.05	0.9	20	112	0
2,4,6-Tribromophenol	0.127	0.198	.05	-55.9*	20	157	0
Phenanthrene-d10	1	1	.05	0	20	126	0
4,6-Dinitro-o-cresol	5	9.924	.05	-98.5*	20	276	0
Hexachlorobenzene	0.228	0.207	.05	9.2	20	115	0
Pentachlorophenol	5	8.051	.05	-61*	20	223	0
Phenanthrene	1.104	0.962	.05	12.9	20	110	0
Anthracene	1.024	0.944	.05	7.8	20	110	0
O-Terphenyl-MS	0.531	0.512	.05	3.6	20	122	0
Fluoranthene	1.146	1.162	.05	-1.4	20	120	0
Pyrene	1.198	1.221	.05	-1.9	20	119	0
4-Terphenyl-d14	0.726	0.711	.05	2.1	20	126	0
Chrysene-d12	1	1	.05	0	20	126	0
Benzo[a]anthracene	1.263	1.173	.05	7.1	20	114	0
Chrysene	1.312	1.257	.05	4.2	20	121	0
Bis(2-ethylhexyl)phthalate	5	7.877	.05	-57.5*	20	206	0
Perylene-d12	1	1	.05	0	20	116	0
Benzo[b]fluoranthene	1.265	1.375	.05	-8.7	20	116	0
Benzo[k]fluoranthene	1.252	1.414	.05	-12.9	20	116	0
Benzo[a]pyrene	1.044	1.281	.05	-22.7*	20	125	0
Indeno[1,2,3-cd]pyrene	0.926	0.891	.05	3.8	20	100	0
Dibenzo[a,h]anthracene	0.989	1.131	.05	-14.4	20	115	0
Benzo[g,h,i]perylene	1.108	1.123	.05	-1.4	20	107	0



^{*} Value outside of QC limits.

Calibration Verification Summary Form 7 Semivolatiles

Client : Paradigm Environmental Services Lab Number : L2228058
Project Name : 222457 Project Number : 222457

Project Name : 222457 Project Number : 222457
Instrument ID : SV115 Calibration Date : 05/31/22 12:43

 Lab File ID
 : CCV0531B
 Init. Calib. Date(s)
 : 02/04/22

 Sample No
 : WG1644744-3
 Init. Calib. Times
 : 15:41
 18:08

Channel :

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
1,4-Dichlorobenzene-d4	1	1	.05	0	20	78	0
2-Fluorophenol	0.944	1.118	.05	-18.4	20	89	0
Phenol-d6	1.196	1.381	.05	-15.5	20	84	0
Hexachloroethane	0.496	0.556	.05	-12.1	20	84	0
Bis(2-chloroethyl)ether	1.127	1.168	.05	-3.6	20	79	0
n-nitrosodi-n-propylamine	0.764	0.915	.05	-19.8	20	86	0
Nitrobenzene-d5	0.917	1.381	.05	-50.6*	20	106	0
Naphthalene-d8	1	1	.05	0	20	81	0
Naphthalene	1.025	1.036	.05	-1.1	20	82	0
Hexachlorobutadiene	0.189	0.198	.05	-4.8	20	86	0
2-Methylnaphthalene	0.627	0.677	.05	-8	20	86	0
1-Methylnaphthalene	0.607	0.637	.05	-4.9	20	86	0
2-Fluorobiphenyl	0.735	0.744	.05	-1.2	20	89	0
2-Chloronaphthalene	0.649	0.669	.05	-3.1	20	84	0
Acenaphthylene	0.972	1.147	.05	-18	20	90	0
Acenaphthene-d10	1	1	.05	0	20	88	0
Acenaphthene	1.325	1.326	.05	-0.1	20	87	0
Fluorene	1.404	1.508	.05	-7.4	20	90	0
2,4,6-Tribromophenol	0.127	0.222	.05	-74.8*	20	130	0
Phenanthrene-d10	1	1	.05	0	20	89	0
4,6-Dinitro-o-cresol	5	10.361	.05	-107.2*	20	205	0
Hexachlorobenzene	0.228	0.236	.05	-3.5	20	93	0
Pentachlorophenol	5	8.297	.05	-65.9*	20	163	0
Phenanthrene	1.104	1.12	.05	-1.4	20	90	0
Anthracene	1.024	1.133	.05	-10.6	20	93	0
O-Terphenyl-MS	0.531	0.578	.05	-8.9	20	98	0
Fluoranthene	1.146	1.291	.05	-12.7	20	94	0
Pyrene	1.198	1.337	.05	-11.6	20	92	0
4-Terphenyl-d14	0.726	0.797	.05	-9.8	20	100	0
Chrysene-d12	1	1	.05	0	20	91	0
Benzo[a]anthracene	1.263	1.391	.05	-10.1	20	97	0
Chrysene	1.312	1.308	.05	0.3	20	91	0
Bis(2-ethylhexyl)phthalate	5	7.775	.05	-55.5*	20	147	0
Perylene-d12	1	1	.05	0	20	94	0
Benzo[b]fluoranthene	1.265	1.447	.05	-14.4	20	99	0
Benzo[k]fluoranthene	1.252	1.383	.05	-10.5	20	92	0
Benzo[a]pyrene	1.044	1.341	.05	-28.4*	20	106	0
Indeno[1,2,3-cd]pyrene	0.926	1.151	.05	-24.3*	20	105	0
Dibenzo[a,h]anthracene	0.989	1.187	.05	-20	20	98	0
Benzo[g,h,i]perylene	1.108	1.226	.05	-10.6	20	95	0



^{*} Value outside of QC limits.

Appendix C

Validator Qualifications

KENNETH R. APPLIN Geochemist/Data Validator

Ph.D., Geochemistry and Mineralogy, The Pennsylvania State University

M.S., Geochemistry and Mineralogy, The Pennsylvania State University

B.A., Geological Sciences, SUNY at Geneseo, NY

Dr. Applin has over 35 years of experience working with the geochemistry of natural waters. His prior experience includes working as an Assistant Professor of Geology at the University of Missouri-Columbia and as Chief Hydrogeologist and Geochemist with a leading engineering firm in Rochester, NY. In 1993, he established KR Applin and Associates, a small consulting business that focuses on the geochemistry of natural waters, especially as applied to problems involving the contamination of groundwater and surface water.

Dr. Applin is also an experienced analytical data validator and has provided data validation services since 1994 to a variety of clients performing brownfield cleanup projects, hazardous waste remediation, groundwater monitoring at solid waste facilities, and other projects requiring third-party data validation. Dr. Applin has several years of hands-on experience with the laboratory analysis of natural waters and has successfully completed the USEPA Region II certification courses for performing inorganic and organic analytical data validation.

MICHAEL K. PERRY Chemist/Data Validator

B.S. Chemistry, Georgia State University, Atlanta, GA

A.A.S., Chemical Technology, Alfred State College, Alfred, NY

Mr. Perry has over 30 years of experience in the analytical laboratory business. During his early career, he spent several years as a laboratory analyst performing the analysis of soil, water, and air samples for inorganic and organic chemical parameters. During his last 20 years in the environmental laboratory business, he managed and directed two major analytical laboratories in Rochester, NY. His management responsibilities included oversight of the daily operations of the lab, staff training and supervision, the selection, purchase, and maintenance of analytical instruments, the introduction of new laboratory methods, analytical quality assurance and quality control, data acquisition and management, and other business-related activities.

Mr. Perry has an extensive working knowledge of the methods and procedures used for sampling and analyzing both inorganic and organic analytes in soil, water, and air. He is an accomplished laboratory chemist and is familiar with the analytical methods and procedures established under the USEPA Contract Laboratory Protocols (CLP), the NYSDEC Analytical Services Protocols (ASP), and the NYSDOH Environmental Laboratory Approval Program (ELAP).