

TEMPORARY GROUNDWATER DISCHARGE PERMIT APPLICATION Drinking Water and Groundwater Bureau Groundwater Discharge Program



RSA/Rule: RSA 485-A:6, VII; 485:3, X; Env-Wq 402

The TEMPORARY GROUNDWATER DISCHARGE PERMIT is a nonrenewable permit issued under RSA 485-A:13 and Env-Wq 402 for the temporary discharge of nondomestic wastewater including that which has received treatment by best available technology (Examples include groundwater remediation, dewatering projects, pump tests, discharges or treated water to the ground or groundwater, etc.)

SUBMIT:

- ONE SIGNED AND COMPLETED APPLICATION TO THE MUNICIPALITY IN WHICH THE DISCHARGE WILL OCCUR
- ONE SIGNED AND COMPLETED APPLICATION TO NHDES AT THE ADDRESS BELOW
- TO: NHDES/Water Division
 Drinking Water & Groundwater Bureau
 Discharge Permit Coordinator
 P.O. Box 95
 Concord, NH 03302-0095

FOR STATE USE ONLY

Date Received: _____

Site No:

Rivers Coordinator Notified Date:

If you have any questions, please contact the Discharge Permits Coordinator at (603) 271-2858.

CERTIFICATION OF MUNICIPAL NOTIFICATION

In order to meet the requirements of Env-Wq 402, the undersigned certifies that on ______(date), a copy of this completed permit application was delivered to the Town/City Clerk of ______(the town in which the proposed discharge will be located).

Date: <u>12/10/2020</u>

Signed:

Applicant (Landowner)

I. Facility

Name: <u>Coakley Landfill</u> Address: <u>480 Breakfast Hill Road</u> City/Town: <u>North Hampton</u> State: <u>NH</u> Zip: <u>03840</u> Latitude and Longitude of Discharge point(s): <u>43.001914, -70.818054</u>

> dwgbinfo@des.nh.gov or phone (603) 271-2858 PO Box 95, Concord, NH 03302-0095 www.des.nh.gov

Property Tax Map: 21 Lot Number: 32, 33

Name: Coakley Landfill Group (Mr. Peter Britz – Contact) Daytime Telephone: (603) 610-7215 Fax Number: ()	II. Applicant (if you are a contact person for the applicant check this box	< 📕)		
Daytime Telephone: (<u>603) 610-7215</u> Fax Number: ()	Name: Coakley Landfill Group (Mr. Peter Britz – Contact)	-		
Mailing Address: 1 Junkins Avenue City/Town: Portsmouth State: NH Zip: 03801 Email Address (Contact Person): plbritz@cityofportsmouth.com Contact Person Phone Number: As above Fax Number: ()	Daytime Telephone: <u>(603) 610-7215</u> Fax Number: ()		
City/Town: Portsmouth State: NH Zip: 03801 Email Address (Contact Person): plbritz@cityofportsmouth.com Contact Person Phone Number: As above Fax Number: () III. Facility Owner (complete only if different from Applicant) Owner Name: Owner Name: City/Town: City/Town: Email Address (Contact Person): Contact Person Phone Number: (Mailing Address: <u>1 Junkins Avenue</u>			
Email Address (Contact Person): plbritz@cityofportsmouth.com Contact Person Phone Number: As above Fax Number: plbritz@cityofportsmouth.com III. Facility Owner (complete only if different from Applicant) Owner	City/Town: Portsmouth State: NH Zip: 03801			
Contact Person Phone Number: <u>As above</u> Fax Number: () III. Facility Owner (complete only if different from Applicant) Owner Name: © Owner © Operator Daytime Telephone: () Mailing Address: State: Zip: City/Town: State: Zip: Email Address (Contact Person): Fax Number: () IV. Property Owner (complete only if different from Applicant or Facility Owner) Name: Daytime Telephone: () Mailing Address: City/Town: State:Zip: City/Town: State:Zip: Email Address (Contact Person): State:Zip:	Email Address (Contact Person): plusted:plust Plusted::plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted:plusted: Plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plusted::plus	<u>1</u>		
III. Facility Owner (complete only if different from Applicant) Owner Name: Owner Owner<	Contact Person Phone Number: <u>As above</u> Fax Number: ()		
Owner Name: Owner Operator Daytime Telephone:	III. Facility Owner (complete only if different from Applicant)			
Daytime Telephone: ()	Owner Name:		Owner	Operator
Mailing Address:	Daytime Telephone: ()			
City/Town: State: Zip: Email Address (Contact Person): Fax Number:	Mailing Address:			
Email Address (Contact Person):	City/Town:	State:	Zip:_	
Contact Person Phone Number:() Fax Number: () IV. Property Owner (complete only if different from Applicant or Facility Owner) Name:	Email Address (Contact Person):			
IV. Property Owner (complete only if different from Applicant or Facility Owner) Name:	Contact Person Phone Number:()	Fax Number:	: ()	
Name:	IV. Property Owner (complete only if different from Applicant or Facility	Owner)		
Daytime Telephone: ()	Name:			
Mailing Address:	Daytime Telephone: ()			
City/Town: State: Zip: Email Address (Contact Person):	Mailing Address:			
Email Address (Contact Person):	City/Town:	State:	Zip:	
Contact Person Phone Number:() Fax Number: ()	Email Address (Contact Person):			
	Contact Person Phone Number:()	Fax Number:	: ()	
V Please provide the following information related to the proposed temporary discharge:	V. Please provide the following information related to the proposed term	norary discha	Irge.	

- a. The purpose of the temporary discharge (e.g. groundwater remediation, well rehab or pumping test, construction dewatering, etc.) <u>Groundwater monitoring well redevelopment.</u>
- b. Proposed Discharge Location

Include a clear color copy of a USGS topographic map or equivalent map which depicts the facility or site location, the discharge location and the location of the closest sanitary sewer.

<u>See attached Figure 1. Closest sanitary sewer is believed to be located along Lafayette Rd and approximately 1,650 ft east of the proposed discharge location.</u>

Location of discharge, if different from facility:			
Address:			_
City/Town:	State:	Zip:	_
Property Tax Map:	Lot Number:		
Latitude & Longitude:			

- c. Location of closest sanitary sewer: Along Lafayette Road, approximately 1,900' east of the discharge location
- d. Proposed Discharge Rate
 - Proposed starting date: <u>One time discharge once permit is approved</u> Estimated discharge: _____5____ gpm for ___4.5____ hours total. Estimated number of days discharge will be required: <u>0.5 days</u>

dwgbinfo@des.nh.gov or phone (603) 271-2858 PO Box 95, Concord, NH 03302-0095 www.des.nh.gov e. Proposed Discharge Method

Describe the method and materials used for the temporary discharge, include a description of any erosion control measures used at the point of discharge: 1,300 gallons of groundwater are currently contained in two plastic storage vessels at the Site immediately adjacent to a landfill stormwater detention pond (northwest corner of landfill). The vessels will be drained via a valve located at the bottom of each vessel. A standard garden hose will be attached to the valve and run to the bottom of the pond and terminated in a perforated 5-gallon bucket. Once opened, the water will be emptied at a rate of no more than 5 gallons per minute with the bucket aiding in dispersion of water and reduction of potential for erosion. Once discharged to the stormwater pond, water drains primarily via infiltration.

- VI. Groundwater Contamination Information, Treatment and Discharge Monitoring
 - a. Provide a summary of the most recent groundwater monitoring results, including total VOCs (laboratory results should also be attached to the application) of the source water for the temporary groundwater discharge:

Location	Compound(s) Exceeding Water Quality Standards	Concentrations (ug/L)
MW-6	Manganese	4140

b. Proposed Treatment

Type of treatment proposed (include a description of the wastewater, information on influent and effluent water quality and on sludge or other by-products generated:

No treatment of water is anticipated prior to disposal.

c. Provide a description of the proposed monitoring and sampling program for the water discharged at the site (applicable only if the source water for the discharge is known to contain, or is anticipated to contain, contamination):

The water will be discharged to a location within an existing Groundwater Management Zone (GMZ) with the area monitored routinely for groundwater and surface water quality in accordance with a USEPA-approved Sampling and Analysis Plan for the Coakley Landfill Superfund Site. Water discharge from the basin, if any, is included as part of ongoing stormwater sampling performed at the landfill and specific to water exiting stormwater control basins.

By signing this application the signer certifies that the information contained in or otherwise submitted with this application is true, complete and not misleading to the best of the signer's knowledge and belief.

By signing this application the signer understands that submission of false, incomplete or misleading information is grounds for:

- Denying the application;
- Revoking any application that is granted based on the information; and
- If the signer is acting as, or on behalf of, a listed engineer as defined in Env-C 502.10, debarring the listed engineer from the roster.

By signing the application, the signer and applicant agree to comply with all applicable rules and conditions of this permit and to not discharge to the holding tank(s) until written permission from the department has been received.

to Buch

Signature of Applicant or Contact Date Signature of Facility Owner (if not Applicant) Date Signature of Property Owner (if not Applicant or Facility Owner) Date

12/10/2020



FIGURE 1

2019-09-08 10424.024



ANALYTICAL REPORT

Lab Number:	L2027711
Client:	CES, Inc 415 Lisbon St. 2nd floor Lewiston, ME 04240
ATTN:	Chris Buckman
Phone:	(207) 989-4824
Project Name:	COAKLEY LANDFILL
Project Number:	10424
Report Date:	07/15/20

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



Serial_No:07152017:03

Project Name: COAKLEY LANDFILL Lab Number: Project Number: 10424 Report Date:

Alpha Sample ID L2027711-01

GW-MW-6-DISP

Client ID

Matrix WATER Sample Location GREENLAND, NH Collection Date/Time 06/30/20 10:30

Receive Date 06/30/20

L2027711

07/15/20



Lab Number: L2027711 Report Date: 07/15/20

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.



 Lab Number:
 L2027711

 Report Date:
 07/15/20

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.

Perfluorinated Alkyl Acids by Isotope Dilution

L2027711-01: The sample was re-analyzed due to QC failures in the original analysis. The results of the reanalysis are reported.

L2027711-01: Extracted Internal Standard recoveries were outside the acceptance criteria for individual analytes. Please refer to the surrogate section of the report for details.

L2027711-01 (MeOH): The sample was re-analyzed due to QC failures in the original analysis. The results of the re-analysis are reported.

L2027711-01 (MeOH): Extracted Internal Standard recoveries were outside the acceptance criteria for individual analytes. Please refer to the surrogate section of the report for details.

L2027711-01: The MeOH fraction of the extraction is reported for the following compounds: N-Methyl Perfluorooctane Sulfonamide (NMeFOSA) and N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA) due to better extraction efficiency of the Surrogates (Extracted Internal Standards).

WG1388862-1R, WG1388862-2 and WG1388862-3: The sample was re-analyzed due to QC failures in the original analysis. The results of the re-analysis are reported.

WG1388862-1R, WG1388862-2 and WG1388862-3: Extracted Internal Standard recoveries were outside the acceptance criteria for individual analytes. Please refer to the surrogate section of the report for details. WG1388862-1, WG1388862-2, and WG1388862-3 (MeOH): The sample was re-analyzed due to QC failures in the original analysis. The results of the re-analysis are reported.

WG1388862-1, WG1388862, and WG1388862-3 (MeOH): Extracted Internal Standard recoveries were outside the acceptance criteria for individual analytes. Please refer to the surrogate section of the report for details.

WG1388862-1/-2/-3: The MeOH fraction of the extraction is reported for the following compounds: N-Methyl Perfluorooctane Sulfonamide (NMeFOSA) and N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA), due to better



 Lab Number:
 L2027711

 Report Date:
 07/15/20

Case Narrative (continued)

extraction efficiency of the Surrogates (Extracted Internal Standards).

WG1388862-2/-3: The internal Standard M2PFDA (form 8) is within the method criteria.

The WG1388862-2/-3 LCS/LCSD RPD, associated with L2027711-01, is above the acceptance criteria for n-

ethyl perfluorooctanesulfonamido ethanol (netfose) (36%).

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:

Michelle M. Monig Michelle M. Morris

Title: Technical Director/Representative

Date: 07/15/20



ORGANICS



VOLATILES



			Serial_No	0:07152017:03
Project Name:	COAKLEY LANDFILL		Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20
		SAMPLE RESULTS		
Lab ID:	L2027711-01		Date Collected:	06/30/20 10:30
Client ID:	GW-MW-6-DISP		Date Received:	06/30/20
Sample Location:	GREENLAND, NH		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water			
Analytical Method:	1,8260C			
Analytical Date:	07/08/20 11:35			
Analyst:	PD			

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Wes	tborough Lab						
Methylene chloride	ND		ug/l	3.0	0.68	1	
1,1-Dichloroethane	ND		ug/l	0.75	0.21	1	
Chloroform	0.45	J	ug/l	0.75	0.22	1	
Carbon tetrachloride	0.94		ug/l	0.50	0.13	1	
1,2-Dichloropropane	ND		ug/l	1.8	0.14	1	
Dibromochloromethane	ND		ug/l	0.50	0.15	1	
1,1,2-Trichloroethane	ND		ug/l	0.75	0.14	1	
Tetrachloroethene	ND		ug/l	0.50	0.18	1	
Chlorobenzene	ND		ug/l	0.50	0.18	1	
Trichlorofluoromethane	ND		ug/l	2.5	0.16	1	
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1	
1,1,1-Trichloroethane	ND		ug/l	0.50	0.16	1	
Bromodichloromethane	ND		ug/l	0.50	0.19	1	
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1	
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1	
1,3-Dichloropropene, Total	ND		ug/l	0.50	0.14	1	
1,1-Dichloropropene	ND		ug/l	2.5	0.24	1	
Bromoform	ND		ug/l	2.0	0.25	1	
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1	
Benzene	ND		ug/l	0.50	0.16	1	
Toluene	ND		ug/l	0.75	0.20	1	
Ethylbenzene	ND		ug/l	0.50	0.17	1	
Chloromethane	ND		ug/l	2.5	0.20	1	
Bromomethane	ND		ug/l	1.0	0.26	1	
Vinyl chloride	ND		ug/l	1.0	0.07	1	
Chloroethane	ND		ug/l	1.0	0.13	1	
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1	
trans-1,2-Dichloroethene	ND		ug/l	0.75	0.16	1	



					S	Serial_No	0:07152017:03	
Project Name:	COAKLEY LANDFILL				Lab Nu	mber:	L2027711	
Project Number:	10424				Report	Date:	07/15/20	
•		SAMP	LE RESULT	5	•		01/10/20	
Lab ID:	L2027711-01				Date Col	lected:	06/30/20 10:30	
Client ID:	GW-MW-6-DISP				Date Red	ceived:	06/30/20	
Sample Location:	GREENLAND, NH				Field Pre	p:	Not Specified	
Sample Depth:			•					
Parameter		Result	Qualifier	Units	RL	MDL	Dilution Factor	_
Volatile Organics b	by GC/MS - Westborough	Lab						
1,2-Dichloroethene, Tota	ıl	ND		ug/l	0.50	0.16	1	
Trichloroethene		ND		ug/l	0.50	0.18	1	
1,2-Dichlorobenzene		ND		ug/l	2.5	0.18	1	
1,3-Dichlorobenzene		ND		ug/l	2.5	0.19	1	
1,4-Dichlorobenzene		ND		ug/l	2.5	0.19	1	
Methyl tert butyl ether		0.92	J	ug/l	1.0	0.17	1	
p/m-Xylene		ND		ug/l	1.0	0.33	1	
o-Xylene		ND		ug/l	1.0	0.39	1	
Xylenes, Total		ND		ug/l	1.0	0.33	1	
cis-1,2-Dichloroethene		ND		ug/l	0.50	0.19	1	
Dibromomethane		ND		ug/l	5.0	0.36	1	
1,2,3-Trichloropropane		ND		ug/l	5.0	0.18	1	
Styrene		ND		ug/l	1.0	0.36	1	
Dichlorodifluoromethane		ND		ug/l	5.0	0.24	1	
Acetone		ND		ug/l	5.0	1.5	1	
Carbon disulfide		ND		ug/l	5.0	0.30	1	
2-Butanone		ND		ug/l	5.0	1.9	1	
4-Methyl-2-pentanone		ND		ug/l	5.0	0.42	1	
2-Hexanone		ND		ug/l	5.0	0.52	1	
Bromochloromethane		ND		ug/l	2.5	0.15	1	
Tetrahydrofuran		ND		ug/l	5.0	0.52	1	
2,2-Dichloropropane		ND		ug/l	2.5	0.20	1	
1,2-Dibromoethane		ND		ug/l	2.0	0.19	1	
1,1,1,2-Tetrachloroethan	e	ND		ug/l	0.50	0.16	1	
Bromobenzene		ND		ug/l	2.5	0.15	1	
n-Butylbenzene		ND		ug/l	0.50	0.19	1	
sec-Butylbenzene		ND		ug/l	0.50	0.18	1	
tert-Butylbenzene		ND		ug/l	2.5	0.20	1	
o-Chlorotoluene		ND		ug/l	2.5	0.22	1	
p-Chlorotoluene		ND		ug/l	2.5	0.18	1	
1,2-Dibromo-3-chloropro	pane	ND		ug/l	2.5	0.35	1	
Hexachlorobutadiene		ND		ug/l	0.50	0.22	1	
Isopropylbenzene		ND		ug/l	0.50	0.19	1	
p-lsopropyltoluene		ND		ug/l	0.50	0.19	1	
Naphthalene		ND		ug/l	2.5	0.22	1	
n-Propylbenzene		ND		ug/l	0.50	0.17	1	
1,2,3-Trichlorobenzene		ND		ua/l	2.5	0.23	1	



			Serial_No:07152017:03				0:07152017:03
Project Name:	COAKLEY LANDFILL				Lab Nu	mber:	L2027711
Project Number:	10424				Report	Date:	07/15/20
		SAMP	LE RESULTS	5			
Lab ID:	L2027711-01				Date Coll	ected:	06/30/20 10:30
Client ID:	GW-MW-6-DISP				Date Rec	eived:	06/30/20
Sample Location:	GREENLAND, NH				Field Pre	o:	Not Specified
Sample Depth:							
Parameter		Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics b	oy GC/MS - Westborough	Lab					
1,2,4-Trichlorobenzene		ND		ug/l	2.5	0.22	1
1,3,5-Trimethylbenzene		ND		ug/l	2.5	0.22	1

ug/l

ug/l

ug/l

ug/l

ug/l

ug/l

ug/l

ug/l

2.0

2.5

2.5

2.0

10

2.0

2.0

250

0.14

0.19

0.16

0.42

1.4

0.18

0.28

61.

ND

ND

ND

ND

36

ND

ND

ND

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	109	70-130	
Toluene-d8	103	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	102	70-130	



1

1

1

1

1

1

1

1

1,3,5-Trichlorobenzene

1,2,4-Trimethylbenzene

Ethyl ether

Isopropyl Ether

1,4-Dioxane

Tert-Butyl Alcohol

Ethyl-Tert-Butyl-Ether

Tertiary-Amyl Methyl Ether

Project Number: 10424

 Lab Number:
 L2027711

 Report Date:
 07/15/20

Method Blank Analysis Batch Quality Control

Analytical Method:1,8260CAnalytical Date:07/08/20 08:47Analyst:PD

Parameter	Result	Qualifier Units	RL	MDL
/olatile Organics by GC/MS	- Westborough Lab	o for sample(s): 01	Batch:	WG1390024-5
Methylene chloride	ND	ug/l	3.0	0.68
1,1-Dichloroethane	ND	ug/l	0.75	0.21
Chloroform	ND	ug/l	0.75	0.22
Carbon tetrachloride	ND	ug/l	0.50	0.13
1,2-Dichloropropane	ND	ug/l	1.8	0.14
Dibromochloromethane	ND	ug/l	0.50	0.15
1,1,2-Trichloroethane	ND	ug/l	0.75	0.14
Tetrachloroethene	ND	ug/l	0.50	0.18
Chlorobenzene	ND	ug/l	0.50	0.18
Trichlorofluoromethane	ND	ug/l	2.5	0.16
1,2-Dichloroethane	ND	ug/l	0.50	0.13
1,1,1-Trichloroethane	ND	ug/l	0.50	0.16
Bromodichloromethane	ND	ug/l	0.50	0.19
trans-1,3-Dichloropropene	ND	ug/l	0.50	0.16
cis-1,3-Dichloropropene	ND	ug/l	0.50	0.14
1,3-Dichloropropene, Total	ND	ug/l	0.50	0.14
1,1-Dichloropropene	ND	ug/l	2.5	0.24
Bromoform	ND	ug/l	2.0	0.25
1,1,2,2-Tetrachloroethane	ND	ug/l	0.50	0.17
Benzene	ND	ug/l	0.50	0.16
Toluene	ND	ug/l	0.75	0.20
Ethylbenzene	ND	ug/l	0.50	0.17
Chloromethane	ND	ug/l	2.5	0.20
Bromomethane	ND	ug/l	1.0	0.26
Vinyl chloride	ND	ug/l	1.0	0.07
Chloroethane	ND	ug/l	1.0	0.13
1,1-Dichloroethene	ND	ug/l	0.50	0.17
trans-1,2-Dichloroethene	ND	ug/l	0.75	0.16
1,2-Dichloroethene, Total	ND	ug/l	0.50	0.16



Project Number: 10424

 Lab Number:
 L2027711

 Report Date:
 07/15/20

Method Blank Analysis Batch Quality Control

Analytical Method:1,8260CAnalytical Date:07/08/20 08:47Analyst:PD

Parameter	Result	Qualifier Units	RL	MDL
Volatile Organics by GC/MS -	Westborough Lab	for sample(s):	01 Batch:	WG1390024-5
Trichloroethene	ND	ug/l	0.50	0.18
1,2-Dichlorobenzene	ND	ug/l	2.5	0.18
1,3-Dichlorobenzene	ND	ug/l	2.5	0.19
1,4-Dichlorobenzene	ND	ug/l	2.5	0.19
Methyl tert butyl ether	ND	ug/l	1.0	0.17
p/m-Xylene	ND	ug/l	1.0	0.33
o-Xylene	ND	ug/l	1.0	0.39
Xylenes, Total	ND	ug/l	1.0	0.33
cis-1,2-Dichloroethene	ND	ug/l	0.50	0.19
Dibromomethane	ND	ug/l	5.0	0.36
1,2,3-Trichloropropane	ND	ug/l	5.0	0.18
Styrene	ND	ug/l	1.0	0.36
Dichlorodifluoromethane	ND	ug/l	5.0	0.24
Acetone	ND	ug/l	5.0	1.5
Carbon disulfide	ND	ug/l	5.0	0.30
2-Butanone	ND	ug/l	5.0	1.9
4-Methyl-2-pentanone	ND	ug/l	5.0	0.42
2-Hexanone	ND	ug/l	5.0	0.52
Bromochloromethane	ND	ug/l	2.5	0.15
Tetrahydrofuran	ND	ug/l	5.0	0.52
2,2-Dichloropropane	ND	ug/l	2.5	0.20
1,2-Dibromoethane	ND	ug/l	2.0	0.19
1,1,1,2-Tetrachloroethane	ND	ug/l	0.50	0.16
Bromobenzene	ND	ug/l	2.5	0.15
n-Butylbenzene	ND	ug/l	0.50	0.19
sec-Butylbenzene	ND	ug/l	0.50	0.18
tert-Butylbenzene	ND	ug/l	2.5	0.20
o-Chlorotoluene	ND	ug/l	2.5	0.22
p-Chlorotoluene	ND	ug/l	2.5	0.18



Project Number:

10424

Lab Number: **Report Date:**

L2027711 07/15/20

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 07/08/20 08:47 Analyst: PD

Parameter	Result C	Qualifier Units	RL	MDL
Volatile Organics by GC/MS - West	borough Lab f	or sample(s): 01	Batch:	WG1390024-5
1,2-Dibromo-3-chloropropane	ND	ug/l	2.5	0.35
Hexachlorobutadiene	ND	ug/l	0.50	0.22
Isopropylbenzene	ND	ug/l	0.50	0.19
p-Isopropyltoluene	ND	ug/l	0.50	0.19
Naphthalene	ND	ug/l	2.5	0.22
n-Propylbenzene	ND	ug/l	0.50	0.17
1,2,3-Trichlorobenzene	ND	ug/l	2.5	0.23
1,2,4-Trichlorobenzene	ND	ug/l	2.5	0.22
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.22
1,3,5-Trichlorobenzene	ND	ug/l	2.0	0.14
1,2,4-Trimethylbenzene	ND	ug/l	2.5	0.19
Ethyl ether	ND	ug/l	2.5	0.16
Isopropyl Ether	ND	ug/l	2.0	0.42
Tert-Butyl Alcohol	ND	ug/l	10	1.4
Ethyl-Tert-Butyl-Ether	ND	ug/l	2.0	0.18
Tertiary-Amyl Methyl Ether	ND	ug/l	2.0	0.28
1,4-Dioxane	ND	ug/l	250	61.

Surrogate	%Recovery	A Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	109		70-130
Toluene-d8	105		70-130
4-Bromofluorobenzene	98		70-130
Dibromofluoromethane	92		70-130



Project Number: 10424 Lab Number: L2027711

	LCS		LCSD		%Recovery		F	PD
Parameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual Li	mits
Volatile Organics by GC/MS	- Westborough Lab Associated	l sample(s): 0'	1 Batch: WG	31390024-3	WG1390024-4			
Methylene chloride	83		85		70-130	2		20
1,1-Dichloroethane	94		97		70-130	3		20
Chloroform	85		86		70-130	1		20
Carbon tetrachloride	89		90		63-132	1		20
1,2-Dichloropropane	84		90		70-130	7		20
Dibromochloromethane	83		87		63-130	5		20
1,1,2-Trichloroethane	86		89		70-130	3		20
Tetrachloroethene	91		96		70-130	5		20
Chlorobenzene	93		95		75-130	2		25
Trichlorofluoromethane	94		95		62-150	1		20
1,2-Dichloroethane	88		92		70-130	4		20
1,1,1-Trichloroethane	85		89		67-130	5		20
Bromodichloromethane	78		80		67-130	3		20
trans-1,3-Dichloropropene	83		84		70-130	1		20
cis-1,3-Dichloropropene	79		82		70-130	4		20
1,1-Dichloropropene	86		90		70-130	5		20
Bromoform	81		82		54-136	1		20
1,1,2,2-Tetrachloroethane	83		88		67-130	6		20
Benzene	85		91		70-130	7		25
Toluene	91		95		70-130	4		25
Ethylbenzene	93		96		70-130	3		20
Chloromethane	84		85		64-130	1		20
Bromomethane	93		97		39-139	4		20



Project Name: COAKLEY LANDFILL

Project Number: 10424 Lab Number: L2027711

	LCS		LCSD		%Recovery			RPD	
Parameter	%Recovery	Qual	%Recove	ery Qual	Limits	RPD	Qual	Limits	
Volatile Organics by GC/MS - Westborough L	ab Associated	I sample(s): 0	1 Batch:	WG1390024-3	WG1390024-4				
Vinyl chloride	81		85		55-140	5		20	
Chloroethane	100		99		55-138	1		20	
1,1-Dichloroethene	84		93		61-145	10		25	
trans-1,2-Dichloroethene	88		90		70-130	2		20	
Trichloroethene	81		83		70-130	2		25	
1,2-Dichlorobenzene	98		98		70-130	0		20	
1,3-Dichlorobenzene	98		98		70-130	0		20	
1,4-Dichlorobenzene	96		96		70-130	0		20	
Methyl tert butyl ether	79		82		63-130	4		20	
p/m-Xylene	95		95		70-130	0		20	
o-Xylene	95		100		70-130	5		20	
cis-1,2-Dichloroethene	86		90		70-130	5		20	
Dibromomethane	82		86		70-130	5		20	
1,2,3-Trichloropropane	88		92		64-130	4		20	
Styrene	90		95		70-130	5		20	
Dichlorodifluoromethane	53		55		36-147	4		20	
Acetone	86		88		58-148	2		20	
Carbon disulfide	85		89		51-130	5		20	
2-Butanone	85		93		63-138	9		20	
4-Methyl-2-pentanone	71		77		59-130	8		20	
2-Hexanone	78		83		57-130	6		20	
Bromochloromethane	88		91		70-130	3		20	
Tetrahydrofuran	100		120		58-130	18		20	



Project Number: 10424 Lab Number: L2027711

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	RP Qual Lim	D hits
Volatile Organics by GC/MS - Westborough I	_ab Associated	sample(s): (1 Batch: WG	1390024-3	WG1390024-4			
2,2-Dichloropropane	90		90		63-133	0	2	0
1,2-Dibromoethane	88		89		70-130	1	2	0
1,1,1,2-Tetrachloroethane	92		91		64-130	1	2	0
Bromobenzene	94		96		70-130	2	2	0
n-Butylbenzene	100		100		53-136	0	2	0
sec-Butylbenzene	99		99		70-130	0	2	0
tert-Butylbenzene	82		82		70-130	0	2	0
o-Chlorotoluene	97		96		70-130	1	2	0
p-Chlorotoluene	96		96		70-130	0	2	0
1,2-Dibromo-3-chloropropane	73		79		41-144	8	2	0
Hexachlorobutadiene	92		98		63-130	6	2	0
Isopropylbenzene	96		97		70-130	1	2	0
p-Isopropyltoluene	97		98		70-130	1	2	0
Naphthalene	76		81		70-130	6	2	0
n-Propylbenzene	97		98		69-130	1	2	0
1,2,3-Trichlorobenzene	85		89		70-130	5	2	0
1,2,4-Trichlorobenzene	86		91		70-130	6	2	0
1,3,5-Trimethylbenzene	97		98		64-130	1	2	0
1,3,5-Trichlorobenzene	94		96		70-130	2	2	0
1,2,4-Trimethylbenzene	95		96		70-130	1	2	0
Ethyl ether	86		91		59-134	6	2	0
Isopropyl Ether	89		91		70-130	2	2	0
Tert-Butyl Alcohol	70		74		70-130	6	2	0



Project Name: COAKLEY LANDFILL

Project Number: 10424

 Lab Number:
 L2027711

 Report Date:
 07/15/20

	LCS		LCSD		%Recovery			RPD	
Parameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual	Limits	
Volatile Organics by GC/MS - Westborough I	Lab Associated	sample(s): C	1 Batch: WG	1390024-3	WG1390024-4				
Ethyl-Tert-Butyl-Ether	83		87		70-130	5		20	
Tertiary-Amyl Methyl Ether	76		80		66-130	5		20	
1,4-Dioxane	86		88		56-162	2		20	

	LCS	LCSD	Acceptance
Surrogate	%Recovery Qu	al %Recovery Qual	Criteria
1,2-Dichloroethane-d4	97	105	70-130
Toluene-d8	102	104	70-130
4-Bromofluorobenzene	99	99	70-130
Dibromofluoromethane	98	99	70-130



SEMIVOLATILES



			Serial_No:	07152017:03
Project Name:	COAKLEY LANDFILL		Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20
		SAMPLE RESULTS		
Lab ID:	L2027711-01		Date Collected:	06/30/20 10:30
Client ID:	GW-MW-6-DISP		Date Received:	06/30/20
Sample Location:	GREENLAND, NH		Field Prep:	Not Specified
Sample Depth:				
Matrix:	Water		Extraction Method:	EPA 3510C
Analytical Method:	1,8270D-SIM		Extraction Date:	07/01/20 08:00
Analytical Date:	07/02/20 23:31			
Analyst:	PS			
-				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1,4 Dioxane by 8270D-SIM - Mansfield Lab						
1,4-Dioxane	0.141	J	ug/l	0.150	0.0339	1
Surrogate			% Recovery	Qualifier	Acce Cr	ptance iteria
1,4-Dioxane-d8			48		1	5-110



			Serial_No	:07152017:03
Project Name:	COAKLEY LANDFILL	-	Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20
			SAMPLE RESULTS	
Lab ID:	L2027711-01	R	Date Collected:	06/30/20 10:30
Client ID:	GW-MW-6-DISP		Date Received:	06/30/20
Sample Location:	GREENLAND, NH		Field Prep:	Not Specified
Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	Water 134,LCMSMS-ID 07/14/20 06:40 RS		Extraction Method Extraction Date:	: ALPHA 23528 07/05/20 11:00

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Perfluorinated Alkyl Acids by Isotope Dilution	n - Mansfield	d Lab				
Perfluorobutanoic Acid (PFBA)	2.17		ng/l	1.83	0.374	1
Perfluoropentanoic Acid (PFPeA)	3.52		ng/l	1.83	0.363	1
Perfluorobutanesulfonic Acid (PFBS)	2.50		ng/l	1.83	0.218	1
Perfluorohexanoic Acid (PFHxA)	12.1		ng/l	1.83	0.300	1
Perfluoroheptanoic Acid (PFHpA)	18.4		ng/l	1.83	0.206	1
Perfluorohexanesulfonic Acid (PFHxS)	1.27	J	ng/l	1.83	0.344	1
Perfluorooctanoic Acid (PFOA)	10.3		ng/l	1.83	0.216	1
1H,1H,2H,2H-Perfluorooctanesulfonic Acid (6:2FTS)	292		ng/l	1.83	1.22	1
Perfluoroheptanesulfonic Acid (PFHpS)	ND		ng/l	1.83	0.630	1
Perfluorononanoic Acid (PFNA)	0.432	J	ng/l	1.83	0.286	1
Perfluorooctanesulfonic Acid (PFOS)	5.91		ng/l	1.83	0.462	1
Perfluorodecanoic Acid (PFDA)	0.619	J	ng/l	1.83	0.278	1
1H,1H,2H,2H-Perfluorodecanesulfonic Acid (8:2FTS)	ND		ng/l	1.83	1.11	1
N-Methyl Perfluorooctanesulfonamidoacetic Acid (NMeFOSAA)	ND		ng/l	1.83	0.593	1
Perfluoroundecanoic Acid (PFUnA)	0.458	J	ng/l	1.83	0.238	1
Perfluorodecanesulfonic Acid (PFDS)	ND		ng/l	1.83	0.897	1
Perfluorooctanesulfonamide (FOSA)	ND		ng/l	1.83	0.531	1
N-Ethyl Perfluorooctanesulfonamidoacetic Acid (NEtFOSAA)	ND		ng/l	1.83	0.736	1
Perfluorododecanoic Acid (PFDoA)	ND		ng/l	1.83	0.341	1
Perfluorotridecanoic Acid (PFTrDA)	ND		ng/l	1.83	0.300	1
Perfluorotetradecanoic Acid (PFTA)	ND		ng/l	1.83	0.227	1
Perfluorohexadecanoic Acid (PFHxDA)	ND		ng/l	3.66	1.14	1
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)	ND		ng/l	45.8	20.3	1
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)	ND		ng/l	45.8	20.6	1



Parameter			Result	Qualifier	Units	RL	MDL	Dilution Factor
Sample Depth:								
Sample Location:	GREENLAND, NH					Field Pre	p:	Not Specified
Client ID:	GW-MW-6-DISP					Date Rec	eived:	06/30/20
Lab ID:	L2027711-01	R				Date Coll	ected:	06/30/20 10:30
			SAMPLE	RESULIS				
Project Number:	10424					Report	Date:	07/15/20
r roject Name.	COARLET LANDFILL						inder.	
Project Name		1				Lah Nu	mher	1 2027711
						S	Serial_No	0:07152017:03

Perfluorinated Alkyl Acids by Isotope Dilution - Mansfield Lab

	2-156
Perfluoro[13C4]Butanoic Acid (MPFBA) 40	
Perfluoro[13C5]Pentanoic Acid (M5PFPEA) 41	16-173
Perfluoro[2,3,4-13C3]Butanesulfonic Acid (M3PFBS) 101	31-159
1H,1H,2H,2H-Perfluoro[1,2-13C2]Hexanesulfonic Acid (M2-4:2FTS) 142	1-313
Perfluoro[1,2,3,4,6-13C5]Hexanoic Acid (M5PFHxA) 31	21-145
Perfluoro[1,2,3,4-13C4]Heptanoic Acid (M4PFHpA) 35	30-139
Perfluoro[1,2,3-13C3]Hexanesulfonic Acid (M3PFHxS) 106	47-153
Perfluoro[13C8]Octanoic Acid (M8PFOA) 38	36-149
1H,1H,2H,2H-Perfluoro[1,2-13C2]Octanesulfonic Acid (M2-6:2FTS) 138	1-244
Perfluoro[13C9]Nonanoic Acid (M9PFNA) 41	34-146
Perfluoro[13C8]Octanesulfonic Acid (M8PFOS) 93	42-146
Perfluoro[1,2,3,4,5,6-13C6]Decanoic Acid (M6PFDA) 44	38-144
1H,1H,2H,2H-Perfluoro[1,2-13C2]Decanesulfonic Acid (M2-8:2FTS) 88	7-170
N-Deuteriomethylperfluoro-1-octanesulfonamidoacetic Acid (d3-NMeFOSAA) 48	1-181
Perfluoro[1,2,3,4,5,6,7-13C7]Undecanoic Acid (M7-PFUDA) 48	40-144
Perfluoro[13C8]Octanesulfonamide (M8FOSA) 16	1-87
N-Deuterioethylperfluoro-1-octanesulfonamidoacetic Acid (d5-NEtFOSAA) 57	23-146
Perfluoro[1,2-13C2]Dodecanoic Acid (MPFDOA) 43	24-161
Perfluoro[1,2-13C2]Tetradecanoic Acid (M2PFTEDA) 45	33-143
Perfluoro[13C2]Hexadecanoic Acid (M2PFHxDA) 70	50-150
2-(N-Methyl-d3-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d7-NMeFOSE) 1	Q 50-150
2-(N-Ethyl-d5-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d9-NEtFOSE) 3	Q 50-150



			Serial_No:07152017:03			
Project Name:	COAKLEY LANDFILL	-		Lab Number:	L2027711	
Project Number:	10424			Report Date:	07/15/20	
			SAMPLE RESULTS			
Lab ID: Client ID: Sample Location:	L2027711-01 GW-MW-6-DISP GREENLAND, NH	R		Date Collected: Date Received: Field Prep:	06/30/20 10:30 06/30/20 Not Specified	
Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	Water 134,LCMSMS-ID 07/15/20 02:23 RS			Extraction Method Extraction Date:	ALPHA 23528 07/05/20 11:00	

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Perfluorinated Alkyl Acids by Isotope Dilution	- Mansfiel	d Lab				
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)	ND		ng/l	18.3	6.74	1
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)	ND		ng/l	18.3	6.08	1
Surrogate (Extracted Internal Standard)			% Recovery	Qualifier	Accej Cri	otance teria
N-Methyl-d3-Perfluoro-1-Octanesulfonamide (d3-NM	eFOSA)		3	Q	5	0-150
N-Ethyl-d5-Perfluoro-1-Octanesulfonamide (d5-NEtFOSA)			2	Q	5	0-150



Serial_No:07152017:03

Project Name: Project Number:	COAKLEY LANDFILL 10424		Lab Number: Report Date:	L2027711 07/15/20
		Method Blank Analysis Batch Quality Control		
Analytical Method: Analytical Date: Analyst:	1,8270D-SIM 07/02/20 19:00 PS		Extraction Method: Extraction Date:	EPA 3510C 07/01/20 08:00

		Quaimer	Units		RL	MDL
,4 Dioxane by 8270D-SIM - Mansfie	eld Lab for	sample(s):	01	Batch:	WG138795	2-1
1,4-Dioxane	ND		ug/l		0.150	0.0339

Surrogate	%Recovery Qualifier	Acceptance Criteria
1,4-Dioxane-d8	43	15-110



Lab Number:

Report Date:

Project Name: COAKLEY LANDFILL

Project Number: 10424

Method Blank Analysis Batch Quality Control

L2027711

07/15/20

Analytical Method:13Analytical Date:07Analyst:JV

134,LCMSMS-ID 07/14/20 11:02 JW Extraction Method: ALPHA 23528 Extraction Date: 07/05/20 11:00

Parameter	Result	Qualifier	Units	RL	MDL	
Perfluorinated Alkyl Acids by Isotope	Dilution -	Mansfield	Lab for sa	ample(s): 01	Batch: W	/G1388862-1 R
Perfluorobutanoic Acid (PFBA)	ND		ng/l	2.00	0.408	
Perfluoropentanoic Acid (PFPeA)	ND		ng/l	2.00	0.396	
Perfluorobutanesulfonic Acid (PFBS)	ND		ng/l	2.00	0.238	
Perfluorohexanoic Acid (PFHxA)	0.348	J	ng/l	2.00	0.328	
Perfluoroheptanoic Acid (PFHpA)	ND		ng/l	2.00	0.225	
Perfluorohexanesulfonic Acid (PFHxS)	ND		ng/l	2.00	0.376	
Perfluorooctanoic Acid (PFOA)	0.264	J	ng/l	2.00	0.236	
1H,1H,2H,2H-Perfluorooctanesulfonic Acid (6:2FTS)	ND		ng/l	2.00	1.33	
Perfluoroheptanesulfonic Acid (PFHpS)	ND		ng/l	2.00	0.688	
Perfluorononanoic Acid (PFNA)	ND		ng/l	2.00	0.312	
Perfluorooctanesulfonic Acid (PFOS)	ND		ng/l	2.00	0.504	
Perfluorodecanoic Acid (PFDA)	ND		ng/l	2.00	0.304	
1H,1H,2H,2H-Perfluorodecanesulfonic Acio (8:2FTS)	I ND		ng/l	2.00	1.21	
N-Methyl Perfluorooctanesulfonamidoacetic Acid (NMeFOSAA)	C ND		ng/l	2.00	0.648	
Perfluoroundecanoic Acid (PFUnA)	ND		ng/l	2.00	0.260	
Perfluorodecanesulfonic Acid (PFDS)	ND		ng/l	2.00	0.980	
Perfluorooctanesulfonamide (FOSA)	ND		ng/l	2.00	0.580	
N-Ethyl Perfluorooctanesulfonamidoacetic Acid (NEtFOSAA)	ND		ng/l	2.00	0.804	
Perfluorododecanoic Acid (PFDoA)	ND		ng/l	2.00	0.372	
Perfluorotridecanoic Acid (PFTrDA)	ND		ng/l	2.00	0.327	
Perfluorotetradecanoic Acid (PFTA)	ND		ng/l	2.00	0.248	
Perfluorohexadecanoic Acid (PFHxDA)	ND		ng/l	4.00	1.24	
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)	ND		ng/l	50.0	22.2	
N-Ethyl Perfluorooctanesulfonamido Ethan (NEtFOSE)	ol ND		ng/l	50.0	22.5	



Project Name:	COAKLEY LANDFILL		Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20
		Method Blank Analysis Batch Quality Control		

Analytical Method:	134,LCMSMS-ID	Extra
Analytical Date:	07/14/20 11:02	Extra
Analyst:	JW	

Extraction Method: ALPHA 23528 Extraction Date: 07/05/20 11:00

Parameter	Result	Qualifier	Units	RL		MDL	
Perfluorinated Alkyl Acids by Isotop	e Dilution - I	Mansfield L	_ab for s	ample(s):	01	Batch:	WG1388862-1 R

Surrogate (Extracted Internal Standard)	%Recovery	/ Qualifier	Acceptance Criteria
Perfluoro[13C4]Butanoic Acid (MPFBA)	60		2-156
Perfluoro[13C5]Pentanoic Acid (M5PFPEA)	68		16-173
Perfluoro[2,3,4-13C3]Butanesulfonic Acid (M3PFBS)	95		31-159
1H,1H,2H,2H-Perfluoro[1,2-13C2]Hexanesulfonic Acid (M2-4:2FTS)	69		1-313
Perfluoro[1,2,3,4,6-13C5]Hexanoic Acid (M5PFHxA)	60		21-145
Perfluoro[1,2,3,4-13C4]Heptanoic Acid (M4PFHpA)	68		30-139
Perfluoro[1,2,3-13C3]Hexanesulfonic Acid (M3PFHxS)	101		47-153
Perfluoro[13C8]Octanoic Acid (M8PFOA)	72		36-149
1H,1H,2H,2H-Perfluoro[1,2-13C2]Octanesulfonic Acid (M2-6:2FTS)	79		1-244
Perfluoro[13C9]Nonanoic Acid (M9PFNA)	74		34-146
Perfluoro[13C8]Octanesulfonic Acid (M8PFOS)	94		42-146
Perfluoro[1,2,3,4,5,6-13C6]Decanoic Acid (M6PFDA)	78		38-144
1H,1H,2H,2H-Perfluoro[1,2-13C2]Decanesulfonic Acid (M2-8:2FTS)	85		7-170
N-Deuteriomethylperfluoro-1-octanesulfonamidoacetic Acid (d3-NMeFOSAA)	55		1-181
Perfluoro[1,2,3,4,5,6,7-13C7]Undecanoic Acid (M7-PFUDA)	86		40-144
Perfluoro[13C8]Octanesulfonamide (M8FOSA)	33		1-87
N-Deuterioethylperfluoro-1-octanesulfonamidoacetic Acid (d5-NEtFOSAA)	57		23-146
Perfluoro[1,2-13C2]Dodecanoic Acid (MPFDOA)	83		24-161
Perfluoro[1,2-13C2]Tetradecanoic Acid (M2PFTEDA)	73		33-143
2,3,3,3-Tetrafluoro-2-[1,1,2,2,3,3,3-Heptafluoropropoxy]-13C3-Propanoic Acid (M3HFPO-DA)	83		50-150
Perfluoro[13C2]Hexadecanoic Acid (M2PFHxDA)	67		50-150
2-(N-Methyl-d3-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d7-NMeFOSE)	22	Q	50-150
2-(N-Ethyl-d5-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d9-NEtFOSE)	21	Q	50-150



Project Name:	COAKLEY LANDFILL		Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20
		Method Blank Analysis Batch Quality Control		

Analytical Method:	134,LCMSMS-ID	Extraction Method:	ALPHA 23528
Analytical Date:	07/14/20 12:08	Extraction Date:	07/05/20 11:00
Analyst:	JW		

Parameter	Result	Qualifier	Units	RL	MDL	
Perfluorinated Alkyl Acids by Isotop	e Dilution -	Mansfield L	ab for s	ample(s): 01	Batch: W	G1388862-1 R
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)	ND		ng/l	20.0	7.36	
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)	ND		ng/l	20.0	6.64	

		A	Acceptance
Surrogate (Extracted Internal Standard)	%Recovery	Qualifier	Criteria
N-Methyl-d3-Perfluoro-1-Octanesulfonamide (d3-NMeFOSA)	1	Q	50-150
N-Ethyl-d5-Perfluoro-1-Octanesulfonamide (d5-NEtFOSA)	1	Q	50-150



Lab Control Sample Analysis

COAKLEY LANDFILL	Batch Quality Control	Lab Number:	L2027711
10424		Report Date:	07/15/20

Parameter	LCS %Recoverv	Qual	LCSD %Recoverv	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
1,4 Dioxane by 8270D-SIM - Mansfield Lab	Associated samp	ole(s): 01	Batch: WG13879	52-2 WG	1387952-3				
1.4-Dioxane	96		102		40-140	6		30	

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,4-Dioxane-d8	44		42		15-110



Project Name:

Project Number:

Project Number: 10424 Lab Number: L2027711

	LCS		LCSD		%Recover	y		RPD
Parameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual	Limits
Perfluorinated Alkyl Acids by Isotope Dilution	- Mansfield Lab	Associated sa	ample(s): 01	Batch: W	VG1388862-2	WG1388862-3		
Perfluorobutanoic Acid (PFBA)	119		116		67-148	7		30
Perfluoropentanoic Acid (PFPeA)	123		119		63-161	4		30
Perfluorobutanesulfonic Acid (PFBS)	122		119		65-157	6		30
Perfluorohexanoic Acid (PFHxA)	123		118		69-168	2		30
Perfluoroheptanoic Acid (PFHpA)	120		117		58-159	5		30
Perfluorohexanesulfonic Acid (PFHxS)	120		115		69-177	8		30
Perfluorooctanoic Acid (PFOA)	123		118		63-159	11		30
1H,1H,2H,2H-Perfluorooctanesulfonic Acid (6:2FTS)	132		134		49-187	4		30
Perfluoroheptanesulfonic Acid (PFHpS)	125		124		61-179	5		30
Perfluorononanoic Acid (PFNA)	122		120		68-171	8		30
Perfluorooctanesulfonic Acid (PFOS)	124		123		52-151	8		30
Perfluorodecanoic Acid (PFDA)	120		116		63-171	4		30
1H,1H,2H,2H-Perfluorodecanesulfonic Acid (8:2FTS)	128		127		56-173	2		30
N-Methyl Perfluorooctanesulfonamidoacetic Acid (NMeFOSAA)	117		116		60-166	12		30
Perfluoroundecanoic Acid (PFUnA)	120		114		60-153	11		30
Perfluorodecanesulfonic Acid (PFDS)	137		131		38-156	16		30
Perfluorooctanesulfonamide (FOSA)	112		115		46-170	15		30
N-Ethyl Perfluorooctanesulfonamidoacetic Acid (NEtFOSAA)	118		114		45-170	3		30
Perfluorododecanoic Acid (PFDoA)	126		118		67-153	6		30
Perfluorotridecanoic Acid (PFTrDA)	119		117		48-158	26		30
Perfluorotetradecanoic Acid (PFTA)	123		118		59-182	13		30
Perfluorohexadecanoic Acid (PFHxDA)	130		137		50-150	0		30



Project Name: COAKLEY LANDFILL

Project Number: 10424

 Lab Number:
 L2027711

 Report Date:
 07/15/20

	LCS		LCSD		%Recover	У		RPD	
Parameter	%Recovery	Qual	%Recovery	' Qual	Limits	RPD	Qual	Limits	
Perfluorinated Alkyl Acids by Isotope Dilution	- Mansfield Lab	Associated sa	ample(s): 01	Batch: \	WG1388862-2	WG1388862-3			
N-Methyl Perfluorooctanesulfonamido Ethanol (NMeFOSE)	114		109		50-150	14		30	
N-Ethyl Perfluorooctanesulfonamido Ethanol (NEtFOSE)	87		92		50-150	36	Q	30	

	LCS		LCSD		Acceptance
Surrogate (Extracted Internal Standard)	%Recovery	Qual	%Recovery	Qual	Criteria
Perfluoro[13C4]Butanoic Acid (MPFBA)	80		84		2-156
Perfluoro[13C5]Pentanoic Acid (M5PFPEA)	85		91		16-173
Perfluoro[2,3,4-13C3]Butanesulfonic Acid (M3PFBS)	90		93		31-159
1H,1H,2H,2H-Perfluoro[1,2-13C2]Hexanesulfonic Acid (M2-4:2FTS)	72		72		1-313
Perfluoro[1,2,3,4,6-13C5]Hexanoic Acid (M5PFHxA)	76		79		21-145
Perfluoro[1,2,3,4-13C4]Heptanoic Acid (M4PFHpA)	81		85		30-139
Perfluoro[1,2,3-13C3]Hexanesulfonic Acid (M3PFHxS)	93		97		47-153
Perfluoro[13C8]Octanoic Acid (M8PFOA)	82		86		36-149
1H,1H,2H,2H-Perfluoro[1,2-13C2]Octanesulfonic Acid (M2-6:2FTS)	79		83		1-244
Perfluoro[13C9]Nonanoic Acid (M9PFNA)	86		87		34-146
Perfluoro[13C8]Octanesulfonic Acid (M8PFOS)	89		89		42-146
Perfluoro[1,2,3,4,5,6-13C6]Decanoic Acid (M6PFDA)	85		84		38-144
1H,1H,2H,2H-Perfluoro[1,2-13C2]Decanesulfonic Acid (M2-8:2FTS)	82		79		7-170
N-Deuteriomethylperfluoro-1-octanesulfonamidoacetic Acid (d3-NMeFOSAA)	72		69		1-181
Perfluoro[1,2,3,4,5,6,7-13C7]Undecanoic Acid (M7-PFUDA)	90		93		40-144
Perfluoro[13C8]Octanesulfonamide (M8FOSA)	32		27		1-87
N-Deuterioethylperfluoro-1-octanesulfonamidoacetic Acid (d5-NEtFOSAA)	80		79		23-146
Perfluoro[1,2-13C2]Dodecanoic Acid (MPFDOA)	88		90		24-161
Perfluoro[1,2-13C2]Tetradecanoic Acid (M2PFTEDA)	75		78		33-143
2,3,3,3-Tetrafluoro-2-[1,1,2,2,3,3,3-Heptafluoropropoxy]-13C3-Propanoic Acid (M3HFPO-DA)	87		83		50-150
Perfluoro[13C2]Hexadecanoic Acid (M2PFHxDA)	72		75		50-150
2-(N-Methyl-d3-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d7-NMeFOSE)	18	Q	14	Q	50-150
2-(N-Ethyl-d5-Perfluoro-1-Octanesulfonamido)ethan-d4-ol (d9-NEtFOSE)	22	Q	16	Q	50-150



Project Name: COAKLEY LANDFILL

Project Number: 10424

 Lab Number:
 L2027711

 Report Date:
 07/15/20

	LCS		LCSD		%Recover	У		RPD	
Parameter	%Recovery	Qual	%Recovery	Qual	Limits	RPD	Qual	Limits	
Perfluorinated Alkyl Acids by Isotope Dilution	- Mansfield Lab	Associated s	ample(s): 01	Batch:	WG1388862-2	WG1388862-3			
N-Methyl Perfluorooctane Sulfonamide (NMeFOSA)	86		97		50-150	12		30	
N-Ethyl Perfluorooctane Sulfonamide (NEtFOSA)	109		103		50-150	6		30	

Surrogate (Extracted Internal Standard)	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria	
N-Methyl-d3-Perfluoro-1-Octanesulfonamide (d3-NMeFOSA)	3	Q	1	Q	50-150	
N-Ethyl-d5-Perfluoro-1-Octanesulfonamide (d5-NEtFOSA)	2	Q	1	Q	50-150	



METALS



Serial_No:07152017:03

Project Name:	COAKLEY LANDFILL	Lab Number:	L2027711
Project Number:	10424	Report Date:	07/15/20
	SAMPLE RE	SULTS	
Lab ID:	L2027711-01	Date Collected:	06/30/20 10:30
Client ID:	GW-MW-6-DISP	Date Received:	06/30/20
Sample Location:	GREENLAND, NH	Field Prep:	Not Specified

Sample Depth:

Matrix:

Water

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Prep Method	Analytical Method	Analyst
Total Metals - Mans	field Lab										
Arsenic, Total	0.2324	J	ug/l	1.000	0.1650	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Barium, Total	15.00		ug/l	1.000	0.1730	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Cadmium, Total	ND		ug/l	0.2000	0.0599	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Chromium, Total	ND		ug/l	1.000	0.1780	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Lead, Total	ND		ug/l	1.000	0.3430	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Manganese, Total	4140		ug/l	1.000	0.4400	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Mercury, Total	ND		ug/l	0.2000	0.0910	1	07/10/20 18:41	07/11/20 13:31	EPA 245.1	3,245.1	AL
Selenium, Total	ND		ug/l	5.000	1.730	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM
Silver, Total	ND		ug/l	0.4000	0.1630	1	07/06/20 19:38	07/07/20 10:34	EPA 3005A	3,200.8	AM



 Lab Number:
 L2027711

 Report Date:
 07/15/20

Method Blank Analysis Batch Quality Control

Parameter	Result Qualifier	Units	RL	MDL	Dilution Factor	Date Prepared	Date Analyzed	Analytical Method	Analyst
Total Metals - Mansfield	Lab for sample(s):	01 Batch	n: WG13	89132-	1				
Arsenic, Total	ND	ug/l	1.000	0.1650	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Barium, Total	ND	ug/l	1.000	0.1730	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Cadmium, Total	ND	ug/l	0.2000	0.0599	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Chromium, Total	ND	ug/l	1.000	0.1780	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Lead, Total	ND	ug/l	1.000	0.3430	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Manganese, Total	ND	ug/l	1.000	0.4400	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Selenium, Total	ND	ug/l	5.000	1.730	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM
Silver, Total	ND	ug/l	0.4000	0.1630	1	07/06/20 19:38	07/07/20 08:46	3,200.8	AM

Prep Information

Digestion Method: EPA 3005A

Parameter Re	esult Qualifier U	Jnits	RL	MDL	Factor	Date Prepared	Date Analyzed	Analytical Method	Analyst
Total Metals - Mansfield Lab	o for sample(s): 01	Batch:	WG13	90880-1	l				
Mercury, Total N	D	ug/l	0.2000	0.0910	1	07/10/20 18:41	07/11/20 13:26	3,245.1	AL

Prep Information

Digestion Method: EPA 245.1



Lab Control Sample Analysis

Batch Quality Control

Project Name: COAKLEY LANDFILL

Project Number: 10424

Lab Number: L2027711 Report Date: 07/15/20

LCS LCSD %Recovery %Recovery Limits %Recovery Qual RPD **RPD** Limits Parameter Qual Qual Total Metals - Mansfield Lab Associated sample(s): 01 Batch: WG1389132-2 85-115 Arsenic, Total 105 --Barium, Total 114 85-115 --Cadmium, Total 108 85-115 --Chromium, Total 104 85-115 --Lead, Total 106 85-115 --Manganese, Total 103 -85-115 -Selenium, Total 106 85-115 --Silver, Total 106 85-115 --Total Metals - Mansfield Lab Associated sample(s): 01 Batch: WG1390880-2

Mercury, Total - 85-115 -



Matrix Spike Analysis

Project Name:	COAKLEY LANDFILL	Batch Quality Control	Lab Number:	L2027711
Project Number:	10424		Report Date:	07/15/20

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Total Metals - Mansfield Lab As	sociated san	nple(s): 01	QC Batch	ID: WG139088	80-3 QC	C Sample	: L2027711-01	Clier	nt ID: GW-M	W-6-DI	SP	
Mercury, Total	ND	5	5.212	104		-	-		70-130	-		20



20

NC

ug/l

Project Name: Project Number:	COAKLEY LANDFILL 10424	La	B Duplicate Analy Batch Quality Control	La Re	b Number port Date	: L2027711 : 07/15/20	
Parameter		Native Sample	Duplicate Sample	Units	RPD	Qual	RPD Limits
Total Metals - Mansfield	Lab Associated sample(s): 01	QC Batch ID: WG139	0880-4 QC Sample: L2	027711-01	Client ID: G	N-MW-6-D	ISP

ND

Mercury, Total

Lab Duplicato Apolycie

ND



Serial_No:07152017:03 Lab Number: L2027711 *Report Date:* 07/15/20

Sample Receipt and Container Information

YES

Were project specific reporting limits specified?

Cooler Information

Cooler	Custody Seal
В	Absent

Container Info	Container Information			Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2027711-01A	Vial HCI preserved	В	NA		3.6	Y	Absent		8260-NH(14)
L2027711-01B	Vial HCl preserved	В	NA		3.6	Y	Absent		8260-NH(14)
L2027711-01C	Vial HCl preserved	В	NA		3.6	Y	Absent		8260-NH(14)
L2027711-01D	Plastic 250ml HNO3 preserved	В	<2	<2	3.6	Y	Absent		BA-2008T-PPB(180),PB-2008T-PPB(180),AG- 2008T-PPB(180),MN-2008T-PPB(180),CD- 2008T-PPB(180),AS-2008T-PPB(180),SE- 2008T-PPB(180),CR-2008T-PPB(180),HG-U- PPB(28)
L2027711-01E	Amber 250ml unpreserved	В	7	7	3.6	Y	Absent		A2-14-DIOXANESIM-PPB(7)
L2027711-01F	Amber 250ml unpreserved	В	7	7	3.6	Y	Absent		A2-14-DIOXANESIM-PPB(7)
L2027711-01G	Plastic 250ml unpreserved	В	NA		3.6	Y	Absent		A2-537-ISOTOPE-36(14)
L2027711-01H	Plastic 250ml unpreserved	В	NA		3.6	Y	Absent		A2-537-ISOTOPE-36(14)



Project Number: 10424

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PFAS PARAMETER SUMMARY

Parameter	Acronym	CAS Number			
PERFLUOROALKYL CARBOXYLIC ACIDS (PFCAs)					
Perfluorooctadecanoic Acid	PFODA	16517-11-6			
Perfluorohexadecanoic Acid	PFHxDA	67905-19-5			
Perfluorotetradecanoic Acid	PFTA	376-06-7			
Perfluorotridecanoic Acid	PFTrDA	72629-94-8			
Perfluorododecanoic Acid	PFDoA	307-55-1			
Perfluoroundecanoic Acid	PFUnA	2058-94-8			
Perfluorodecanoic Acid	PFDA	335-76-2			
Perfluorononanoic Acid	PFNA	375-95-1			
Perfluorooctanoic Acid	PFOA	335-67-1			
Perfluoroheptanoic Acid	PFHpA	375-85-9			
Perfluorohexanoic Acid	PFHxA	307-24-4			
Perfluoropentanoic Acid	PFPeA	2706-90-3			
Perfluorobutanoic Acid	PFBA	375-22-4			
PERFLUOROALKYL SULFONIC ACIDS (PFSAs)					
Perfluorododecanesulfonic Acid	PFDoDS	79780-39-5			
Perfluorodecanesulfonic Acid	PFDS	335-77-3			
Perfluorononanesulfonic Acid	PFNS	68259-12-1			
Perfluorooctanesulfonic Acid	PFOS	1763-23-1			
Perfluoroheptanesulfonic Acid	PFHpS	375-92-8			
Perfluorohexanesulfonic Acid	PFHxS	355-46-4			
Perfluoropentanesulfonic Acid	PFPeS	2706-91-4			
Perfluorobutanesulfonic Acid	PFBS	375-73-5			
FLUOROTELOMERS					
1H,1H,2H,2H-Perfluorododecanesulfonic Acid	10:2FTS	120226-60-0			
1H,1H,2H,2H-Perfluorodecanesulfonic Acid	8:2FTS	39108-34-4			
1H,1H,2H,2H-Perfluorooctanesulfonic Acid	6:2FTS	27619-97-2			
1H,1H,2H,2H-Perfluorohexanesulfonic Acid	4:2FTS	757124-72-4			
PERFLUOROALKANE SULFONAMIDES (FASAs)					
Perfluorooctanesulfonamide	FOSA	754-91-6			
N-Ethyl Perfluorooctane Sulfonamide	NEtFOSA	4151-50-2			
N-Methyl Perfluorooctane Sulfonamide	NMeFOSA	31506-32-8			
PERFLUOROALKANE SULFONYL SUBSTANCES					
N-Ethyl Perfluorooctanesulfonamido Ethanol	NEtFOSE	1691-99-2			
N-Methyl Perfluorooctanesulfonamido Ethanol	NMeFOSE	24448-09-7			
N-Ethyl Perfluorooctanesulfonamidoacetic Acid	NEtFOSAA	2991-50-6			
N-Methyl Perfluorooctanesulfonamidoacetic Acid	NMeFOSAA	2355-31-9			
PER- and POLYFLUOROALKYL ETHER CARBOXYLIC ACIDS					
2,3,3,3-Tetrafluoro-2-[1,1,2,2,3,3,3-Heptafluoropropoxy]-Propanoic Acid	HFPO-DA	13252-13-6			
4,8-Dioxa-3h-Perfluorononanoic Acid	ADONA	919005-14-4			
CHLORO-PERFLUOROALKYL SULFONIC ACIDS					
11-Chloroeicosafluoro-3-Oxaundecane-1-Sulfonic Acid	11CI-PF3OUdS	763051-92-9			
9-Chlorohexadecafluoro-3-Oxanone-1-Sulfonic Acid	9CI-PF3ONS	756426-58-1			



Project Number: 10424

Lab Number: L2027711

Report Date: 07/15/20

GLOSSARY

Acronyms

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.)
EDL	- 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 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.
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.
	- N-Nirosouphenyrannne/Diphenyrannne.
ND	
NP	- Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.
KL	- 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.

Footnotes

Report Format: DU Report with 'J' Qualifiers



Project Number: 10424

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- 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 Waterpreserved 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. 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.
- B 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 concentrations of the analyte at less than ten times (10x) the 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. 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. For NJ-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 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).
- C Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- **D** 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. (DoD and NYSDEC Part 375 PFAS only.)
- 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.
- J 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.
- 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.

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Data Qualifiers

- **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.

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REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.
- 3 Methods for the Determination of Metals in Environmental Samples, Supplement I. EPA/600/R-94/111. May 1994.
- 134 Determination of Selected Perfluorinated Alkyl Acids in Drinking Water by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry (LC/MS/MS) using Isotope Dilution. Alpha SOP 23528.

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.



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 8260C: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.
EPA 8270D: NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine; 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, 1-Methylnaphthalene.
EPA 3C Fixed gases
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.

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.

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.

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