



90 Sargent Drive • New Haven, Connecticut 06511-5966
203-401-6743 or 877-894-5773 (Telephone) • Website: www.rwater.com

Laboratory Certification IDs

CT: PH-0411, NH: 2239, NY: 11867, PA: 68-05519, RI: LAO00339, MA: M-CT004
See website, www.rwalab.com/rwa-lab-certifications, for certified analyte list.

Client: Paul Hennessy
Randolph-Holbrook Joint Water

275 Pond St
Randolph, MA 02368

781-964-9292
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ANALYTICAL REPORT

Project: FEE-RANDOLPHHOLBROOK-23-000009

Report Date: 8/15/2023

This Laboratory is in compliance with the NELAP requirements of procedures used except where indicated .
This report contains results for the analysis tested, under the sampling conditions described on the Chain Of Custody (COC), as received by the laboratory. This report is incomplete unless all pages indicated in the pagination at the bottom of the page are included.

The COC form has been scanned to accompany the analytical report and is an exact copy of the original.
If you are the client above and have any questions concerning this testing, please do not hesitate to contact RWA Client Services at (203) 401-6743 or (877) 894-5773. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.

Technical Representative



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SAMPLE SUMMARY

Sample ID	Customer ID	Collection Date/Time	Receipt Date
300379873	RAW	7/11/2023 1300	7/12/23
300379874	FINISHED	7/11/2023 1300	7/12/23
300379898	FB - RAW	7/11/2023 1300	7/12/23
300379899	FB - FINISHED	7/11/2023 1300	7/12/23

Case Narrative and Comments

All QC passes criteria unless noted in a Comment below.

Samples were received at the appropriate temperature and in accordance with the chain of custody unless noted.

E = Exceeds calibration range **ND** = Non Detect **FB** = Field Blank

RL = Minimum Reporting Level **MDL** = Method Detection Limit

J = The reported result is below RL but greater than the MDL. The reported result is an estimate.

Massachusetts samples for required water quality sampling are included.

Samples and FBs were received in bottles with preservatives Trizma HCL & Trizma base per method requirements.

"FB" added at beginning/end designates "Field Blank" (Field Reagent Blank) for associated Customer ID sample. Field Blank analytes (unless noted) were shown to be less than 1/3 of the RL as per EPA537 or EPA537.1.

Method EPA537 or EPA537.1 Analyte Results, MDL and RL are adjusted to reflect the actual Final (mL) volume used.

<u>Method</u>	<u>CAS#</u>	<u>PFAS Analyte (Acronym)</u>
537, 537.1	1763-23-1	Perfluorooctanesulfonic acid (PFOS)
537, 537.1	335-67-1	Perfluorooctanoic acid (PFOA)
537, 537.1	355-46-4	Perfluorohexanesulfonic acid (PFHxS)
537, 537.1	375-95-1	Perfluorononanoic acid (PFNA)
537, 537.1	375-85-9	Perfluorohepatanoic acid (PFHpA)
537, 537.1	335-76-2	Perfluorodecanoic acid (PFDA)
537, 537.1	375-73-5	Perfluorobutanesulfonic acid (PFBS)
537, 537.1	307-55-1	Perfluorododecanoic acid (PFDoA)
537, 537.1	307-24-4	Perfluorohexanoic acid (PFHxA)
537, 537.1	376-06-7	Perfluorotetradecanoic acid (PFTA)
537, 537.1	72629-94-8	Perfluorotridecanoic acid (PFTTrDA)
537, 537.1	2058-94-8	Perfluoroundecanoic acid (PFUnA)
537, 537.1	2991-50-6	N-ethyl perfluorooctanesulfonamidoacetic acid (NEtFOSAA)
537, 537.1	2355-31-9	N-methyl perfluorooctanesulfonamidoacetic acid (NMeFOSAA)
537.1	763051-92-9	11-chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS)
537.1	756426-58-1	9-chlorohexadecafluoro-3-oxanone-1-sulfonic acid (9Cl-PF3ONS)
537.1	919005-14-4	4,8-dioxa-3H-perfluorononanoic acid (ADONA)
537.1	13252-13-6	Hexafluoropropylene oxide dimer acid (HFPO-DA)

PFAS6 (MassDEP) = sum of PFOS, PFOA, PFHxS, PFNA, PFHpA and PFDA (only include Results at or above the RL)

MassDEP has established a maximum contaminant level (MCL) of 20 ng/L for PFAS6.

Method EPA537 Analyte Results and RL are adjusted to reflect the actual sample Final (mL) volume used.

Sample# 300379899 Field Blank for EPA 537: fails low for d5-NEtFOSAA surrogate recovery; the associated analytes NMeFOSAA and NEtFOSAA are suspect. This surrogate does not affect the data for the PFAS6 MassDEP compounds; the data for PFAS6 is valid.

Sample ID: 300379873

Customer ID: RAW

Collection Date: 07/11/2023 13:00

PWS ID# / LOC ID#: RAW

Project: FEE-RANDOLPHHOLBROOK-23-000009

Analyte	CAS#	Results	MDL	RL	Units	Dilution	Qualifier	Method	Date Time/ Analyzed	Analyst
PFBS	375-73-5	2.13	0.92	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFHxA	307-24-4	3.33	0.70	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFHpA	375-85-9	1.94	0.64	2.06	ng/L	1.03	J	EPA537	7/30/23 2159	CSS/GMA
PFHxS	355-46-4	2.68	0.93	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFOA	335-67-1	4.97	0.74	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFOS	1763-23-1	7.62	0.92	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFNA	375-95-1	1.03	0.65	2.06	ng/L	1.03	J	EPA537	7/30/23 2159	CSS/GMA
PFDA	335-76-2	ND	0.74	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFUnA	2058-94-8	ND	0.94	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
NMeFOSAA	2355-31-9	ND	1.23	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
NEtFOSAA	2991-50-6	ND	1.06	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFDoA	307-55-1	ND	0.85	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFTrDA	72629-94-8	ND	0.93	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFTA	376-06-7	ND	1.00	2.06	ng/L	1.03		EPA537	7/30/23 2159	CSS/GMA
PFAS6 (MassDEP)		15.27	2.06	2.06	ng/L	1.03				
Surrogates					Results	Recovery Limits			Pass/Fail	
13C-PFHxA (SUR) % Recovery					90.90	70 - 130			Pass	
13C-PFDA (SUR) % Recovery					87.70	70 - 130			Pass	
d5-NEtFOSAA (SUR) % Recovery					82.00	70 - 130			Pass	

Sample Extraction Data:

Lab Number (Field ID)	Prep Method	Batch	Final (mL)	Date
300379873	PFAS_537	537_EXT-230717-1	242	07/17/2023

Sample ID: 300379874

Customer ID: FINISHED

Collection Date: 07/11/2023 13:00

PWS ID# / LOC ID#: 4244001 / 10296

Project: FEE-RANDOLPHHOLBROOK-23-000009

Analyte	CAS#	Results	MDL	RL	Units	Dilution	Qualifier	Method	Date Time/ Analyzed	Analyst
PFBS	375-73-5	1.74	0.89	2.0	ng/L	1.0	J	EPA537	7/30/23 2228	CSS/GMA
PFHxA	307-24-4	3.20	0.68	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFHpA	375-85-9	1.88	0.62	2.0	ng/L	1.0	J	EPA537	7/30/23 2228	CSS/GMA
PFHxS	355-46-4	2.61	0.90	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFOA	335-67-1	4.58	0.72	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFOS	1763-23-1	7.30	0.89	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFNA	375-95-1	0.96	0.63	2.0	ng/L	1.0	J	EPA537	7/30/23 2228	CSS/GMA
PFDA	335-76-2	ND	0.72	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFUnA	2058-94-8	ND	0.91	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
NMeFOSAA	2355-31-9	ND	1.19	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
NEtFOSAA	2991-50-6	ND	1.03	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFDoA	307-55-1	ND	0.83	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFTrDA	72629-94-8	ND	0.90	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA
PFTA	376-06-7	ND	0.97	2.0	ng/L	1.0		EPA537	7/30/23 2228	CSS/GMA

PFAS6 (MassDEP) 14.49 2.00 2.00 ng/L 1.0

Surrogates	Results	Recovery Limits	Pass/Fail
13C-PFHxA (SUR) % Recovery	86.20	70 - 130	Pass
13C-PFDA (SUR) % Recovery	82.50	70 - 130	Pass
d5-NEtFOSAA (SUR) % Recovery	77.10	70 - 130	Pass

Sample Extraction Data:

Lab Number (Field ID)	Prep Method	Batch	Final (mL)	Date
300379874	PFAS_537	537_EXT-230717-1	250	07/17/2023

Sample ID: 300379898

Customer ID: FB - RAW

Collection Date: 07/11/2023 13:00

PWS ID# / LOC ID#: FB - RAW

Project: FEE-RANDOLPHHOLBROOK-23-000009

Analyte	CAS#	Results	MDL	RL	Units	Dilution	Qualifier	Method	Date Time/ Analyzed	Analyst
PFBS	375-73-5	ND	0.89	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFHxA	307-24-4	ND	0.68	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFHpA	375-85-9	ND	0.62	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFHxS	355-46-4	ND	0.90	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFOA	335-67-1	ND	0.72	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFOS	1763-23-1	ND	0.89	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFNA	375-95-1	ND	0.63	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFDA	335-76-2	ND	0.72	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFUnA	2058-94-8	ND	0.91	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
NMeFOSAA	2355-31-9	ND	1.19	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
NEtFOSAA	2991-50-6	ND	1.03	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFDoA	307-55-1	ND	0.83	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFTTrDA	72629-94-8	ND	0.90	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFTA	376-06-7	ND	0.97	2.0	ng/L	1.0		EPA537	7/30/23 2214	CSS/GMA
PFAS6 (MassDEP)		ND	2.00	2.00	ng/L	1.0				
Surrogates					Results	Recovery Limits			Pass/Fail	
13C-PFHxA (SUR) % Recovery					100.20	70 - 130			Pass	
13C-PFDA (SUR) % Recovery					88.00	70 - 130			Pass	
d5-NEtFOSAA (SUR) % Recovery					88.70	70 - 130			Pass	

Sample Extraction Data:

Lab Number (Field ID)	Prep Method	Batch	Final (mL)	Date
300379898	PFAS_537	537_EXT-230717-1	250	07/17/2023

Sample ID: 300379899

Customer ID: FB - FINISHED

Collection Date: 07/11/2023 13:00

PWS ID# / LOC ID#: FB - FINISHED

Project: FEE-RANDOLPHHOLBROOK-23-000009

Analyte	CAS#	Results	MDL	RL	Units	Dilution	Qualifier	Method	Date Time/ Analyzed	Analyst
PFBS	375-73-5	ND	0.92	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFHxA	307-24-4	ND	0.70	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFHpA	375-85-9	ND	0.64	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFHxS	355-46-4	ND	0.93	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFOA	335-67-1	ND	0.74	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFOS	1763-23-1	ND	0.92	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFNA	375-95-1	ND	0.65	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFDA	335-76-2	ND	0.74	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFUnA	2058-94-8	ND	0.94	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
NMeFOSAA	2355-31-9	ND	1.23	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
NEtFOSAA	2991-50-6	ND	1.06	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFD _o A	307-55-1	ND	0.85	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFTrDA	72629-94-8	ND	0.93	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFTA	376-06-7	ND	1.00	2.06	ng/L	1.03		EPA537	7/30/23 2243	CSS/GMA
PFAS6 (MassDEP)		ND	2.06	2.06	ng/L	1.03				
Surrogates		Results	Recovery Limits		Pass/Fail					
13C-PFHxA (SUR) % Recovery		85.10	70 - 130		Pass					
13C-PFDA (SUR) % Recovery		73.30	70 - 130		Pass					
d5-NEtFOSAA (SUR) % Recovery		68.30	70 - 130		Fail					

Sample Extraction Data:

Lab Number (Field ID)	Prep Method	Batch	Final (mL)	Date
300379899	PFAS_537	537_EXT-230717-1	242	07/17/2023



90 Sargent Dr, New Haven, CT 06511
 Phone: 203-401-2700
 Fax: 203-401-6799

Chain of Custody Form

Company Name
 RANDOLPH/HOLBROOK
 JOINT WATER

Company Address
 275 POND ST.
 RANDOLPH, MA 012368

Sampler

PO Number

RWA LIMS Number

Date Collected

Time Collected

Sample ID / Sample Location

200379873 7/11/23 1PM ① RAN

874 7/11/23 1PM ② FINISHED

Number of Bottles

5

5

Test Requested, Container, & Preservative

PFAS
 250mL Plastic w/Tris Hydrochloride
 & Tris (Hydroxymethyl) Aminomethane

PFAS 14

4

4

PFAS Poured Field Blank
 250mL Plastic w/Tris Hydrochloride
 & Tris (Hydroxymethyl) Aminomethane

1

1

FB-898

FB-899

State Sample Collected:

CT NY Other (specify) MA

Evidence of Cooling (Circle) Y or N

Cooler Temp °C: 1.9

Container & Preservative Meet Criteria (Circle): Yes/ No

LOT # SLCN1650

LOT # SLCN1665

NOT FOR COMPLIANCE

MA DEP COMPLIANCE

Relinquished By (Signature): Paul Hennessy

Date & Time: 7/11/23 1:15 PM

Date & Time:

Received By (Signature): [Signature]

Date & Time: 7/12/23 8:30 AM

Date & Time:

Comments:

PFAS 14

Fee - RANDOLPH/HOLBROOK-23-000007

South Central Connecticut Regional Water Authority PFAS QA/QC Summary

Extraction Batch QC for: EPA 537.1

MA Lab Cert.#: M-CT004

Extraction Batch Date: 7/30/2023

Sample ID for LFSM/LFSMD: 300381602

Analyte	LFSM %Recovery	LFSMD %Recovery	RPD of LFSM/LFSMD	LRB (MRL is 2)		LFB 10 ng/L %Recovery
	Acceptance Range 70-130%	Acceptance Range 70-130%	Acceptance Limit <30%	Result, ng/L	Meets < 1/3 of MRL criteria?	Acceptance Range 70-130%
PFBS	99.6	99.0	0.6	ND	[Y]	103.8
PFHxA	91.4	96.9	5.8	ND	[Y]	95.6
HFPO-DA	88.3	94.0	6.3	ND	[Y]	93.4
PFHpA	91.6	97.3	6.0	ND	[Y]	94.4
PFHxS	99.8	106.0	6.0	ND	[Y]	101.5
ADONA	90.5	95.8	5.6	ND	[Y]	96.8
PFOA	95.9	100.2	4.4	ND	[Y]	103.5
PFOS	88.0	92.3	4.7	ND	[Y]	95.8
PFNA	93.7	99.2	5.8	ND	[Y]	96.8
9CI-PF3ONS	87.6	89.3	1.9	ND	[Y]	88.8
PFDA	93.6	92.6	1.1	ND	[Y]	94.4
PFUnA	93.0	89.6	3.8	ND	[Y]	96.2
11CI-PF3OUdS	86.9	85.8	1.3	ND	[Y]	87.9
NMeFOSAA	89.4	95.9	7.0	ND	[Y]	82.1
NEtFOSAA	86.2	79.6	7.9	ND	[Y]	94.2
PFDoA	94.0	89.1	5.4	ND	[Y]	90.0
PFTTrDA	93.4	88.7	5.2	ND	[Y]	89.2
PFTA	90.7	86.4	4.9	ND	[Y]	91.4

Surrogate %Recovery

¹³ C ₂ -PFHxA	¹³ C ₃ -HFPO-DA	¹³ C ₂ -PFDA	^d ₅ -NEtFOSAA	
95.50	92.80	90.80	87.80	LFSM
101.80	101.60	93.50	86.00	LFSMD
90.90	89.00	84.70	80.30	LRB
100.00	98.60	90.60	89.20	LFB

Note: The Surrogate %Recovery for Samples and Poured Field Blanks is included on Final Reports.

All Batch QC passes method criteria unless noted in "Comments" section below

Comments: The matrix spike data *is not* from a sample submitted on this Chain of Custody.

LRB = Laboratory Reagent Blank

RPD = relative percent difference

FB = Field Blank

ND = Non-Detect

LFB = Laboratory Fortified Blank

MRL = Method Reporting Level

Results are ng/L (ppt)

LFSM = Laboratory Fortified Sample Matrix

LFSM/LFSMD spike concentration is 20 ng/L, unless noted otherwise

LFSMD = Laboratory Fortified Sample Matrix Duplicate

Surrogate Acceptance Limit is 70 - 130% Recovery