

LAB CHRONICLE

OrderID:	Q3371	OrderDate:	10/16/2025 4:13:00 PM
Client:	Tetra Tech NUS, Inc.	Project:	NWIRP Bethpage 112G08005-WE13
Contact:	Ernie Wu	Location:	D31,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q3371-02	BP-VPB-184-EB-2025 1016	Water			10/16/25			10/16/25
			SVOC-SIMGroup1	8270-Modified		10/20/25	10/20/25	
Q3371-03	VPB184-HYD-202510 10	Water			10/10/25			10/16/25
			SVOC-SIMGroup1	8270-Modified		10/20/25	10/20/25	
Q3371-05	BP-VPB-184-GW-920- 922	Water			10/15/25			10/16/25
			SVOC-SIMGroup1	8270-Modified		10/20/25	10/20/25	
Q3371-06	VPB184-HYD-202510 16	Water			10/16/25			10/16/25
			SVOC-SIMGroup1	8270-Modified		10/20/25	10/20/25	



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Hit Summary Sheet SW-846

SDG No.: Q3371
Client: Tetra Tech NUS, Inc.

Sample ID	Client ID	Parameter	Concentration	C	MDL	LOD	RDL	Units
Client ID : BP-VPB-184-EB-20251016								
Q3371-02	BP-VPB-184-EB-202510 WATER	1,4-Dioxane	0.090	J	0.07	0.22	0.22	ug/L
		Total Svoc :			0.09			
		Total Concentration:			0.09			
Client ID : VPB184-HYD-20251016								
Q3371-06	VPB184-HYD-20251016 WATER	1,4-Dioxane	0.070	J	0.07	0.2	0.2	ug/L
		Total Svoc :			0.07			
		Total Concentration:			0.07			



SAMPLE DATA

Report of Analysis

Client:	Tetra Tech NUS, Inc.		Date Collected:	10/16/25
Project:	NWIRP Bethpage 112G08005-WE13		Date Received:	10/16/25
Client Sample ID:	BP-VPB-184-EB-20251016		SDG No.:	Q3371
Lab Sample ID:	Q3371-02		Matrix:	Water
Analytical Method:	SW8270ESIM	Level : LOW	% Solid:	0
Sample Wt/Vol:	900 mL	Final Vol: 1000 uL	Test:	SVOC-SIMGroup1
Prep Method :		Prep Date: 10/20/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.090	J	1	0.070	0.22	0.22	ug/L	10/20/25 19:45	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.30			30 - 150		76%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.42			30 - 150		105%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.36			55 - 111		91%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.35			53 - 106		87%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.55	*		58 - 132		137%	SPK: 0.4		
INTERNAL STANDARDS		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	6200								
1146-65-2	Naphthalene-d8	16000								
15067-26-2	Acenaphthene-d10	8400								
1517-22-2	Phenanthrene-d10	17100								
1719-03-5	Chrysene-d12	15000								
1520-96-3	Perylene-d12	14500								

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Report of Analysis

Client:	Tetra Tech NUS, Inc.		Date Collected:	10/10/25
Project:	NWIRP Bethpage 112G08005-WE13		Date Received:	10/16/25
Client Sample ID:	VPB184-HYD-20251010		SDG No.:	Q3371
Lab Sample ID:	Q3371-03		Matrix:	Water
Analytical Method:	SW8270ESIM	Level : LOW	% Solid:	0
Sample Wt/Vol:	970 mL	Final Vol: 1000 uL	Test:	SVOC-SIMGroup1
Prep Method :	Prep Date: 10/20/25			

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.21	U	1	0.070	0.21	0.21	ug/L	10/20/25 20:21	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.016	*		30 - 150		4%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.042	*		30 - 150		10%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.34			55 - 111		86%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.34			53 - 106		84%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.55	*		58 - 132		137%	SPK: 0.4		
INTERNAL STANDARDS		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	7050								
1146-65-2	Naphthalene-d8	18600								
15067-26-2	Acenaphthene-d10	9630								
1517-22-2	Phenanthrene-d10	18000								
1719-03-5	Chrysene-d12	14000								
1520-96-3	Perylene-d12	12700								

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J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Report of Analysis

Client: Tetra Tech NUS, Inc.
Project: NWIRP Bethpage 112G08005-WE13
Client Sample ID: BP-VPB-184-GW-920-922
Lab Sample ID: Q3371-05
Analytical Method: SW8270ESIM Level : LOW
Sample Wt/Vol: 790 mL Final Vol: 1000 uL
Prep Method : Prep Date: 10/20/25

Date Collected: 10/15/25
Date Received: 10/16/25
SDG No.: Q3371
Matrix: Water
% Solid: 0
Test: SVOC-SIMGroup1

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.25	U	1	0.080	0.25	0.25	ug/L	10/20/25 20:58	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.29			30 - 150		73%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.41			30 - 150		101%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.35			55 - 111		88%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.33			53 - 106		83%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.55	*		58 - 132		137%	SPK: 0.4		
INTERNAL STANDARDS										
		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	7240								
1146-65-2	Naphthalene-d8	19100								
15067-26-2	Acenaphthene-d10	10100								
1517-22-2	Phenanthrene-d10	20200								
1719-03-5	Chrysene-d12	17500								
1520-96-3	Perylene-d12	17400								

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J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

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D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Report of Analysis

Client:	Tetra Tech NUS, Inc.		Date Collected:	10/16/25
Project:	NWIRP Bethpage 112G08005-WE13		Date Received:	10/16/25
Client Sample ID:	VPB184-HYD-20251016		SDG No.:	Q3371
Lab Sample ID:	Q3371-06		Matrix:	Water
Analytical Method:	SW8270ESIM	Level : LOW	% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 1000 uL	Test:	SVOC-SIMGroup1
Prep Method :		Prep Date: 10/20/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.070	J	1	0.070	0.20	0.20	ug/L	10/20/25 21:34	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.014	*		30 - 150		4%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.043	*		30 - 150		11%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.31			55 - 111		77%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.32			53 - 106		79%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.52			58 - 132		131%	SPK: 0.4		
INTERNAL STANDARDS		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	7490								
1146-65-2	Naphthalene-d8	19900								
15067-26-2	Acenaphthene-d10	10200								
1517-22-2	Phenanthrene-d10	19700								
1719-03-5	Chrysene-d12	16600								
1520-96-3	Perylene-d12	14800								

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

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Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

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D = Dilution

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A = Aldol-Condensation Reaction Products



QC SUMMARY

Surrogate Summary

SW-846

SDG No.: Q3371

Client: Tetra Tech NUS, Inc.

Analytical Method: 8270-Modified

Lab Sample ID	Client ID	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
							Low	High
PB170167BL	PB170167BL	2-Methylnaphthalene-d10	0.4	0.34	85		30	150
		Fluoranthene-d10	0.4	0.33	83		30	150
		Nitrobenzene-d5	0.4	0.36	90		55	111
		2-Fluorobiphenyl	0.4	0.39	98		53	106
		Terphenyl-d14	0.4	0.45	113		58	132
PB170167BS	PB170167BS	2-Methylnaphthalene-d10	0.4	0.40	99		30	150
		Fluoranthene-d10	0.4	0.34	85		30	150
		Nitrobenzene-d5	0.4	0.42	104		55	111
		2-Fluorobiphenyl	0.4	0.44	111	*	53	106
		Terphenyl-d14	0.4	0.51	127		58	132
PB170167BSD	PB170167BSD	2-Methylnaphthalene-d10	0.4	0.36	90		30	150
		Fluoranthene-d10	0.4	0.31	78		30	150
		Nitrobenzene-d5	0.4	0.38	95		55	111
		2-Fluorobiphenyl	0.4	0.42	105		53	106
		Terphenyl-d14	0.4	0.47	117		58	132
Q3371-02	BP-VPB-184-EB-20251016	2-Methylnaphthalene-d10	0.4	0.30	76		30	150
		Fluoranthene-d10	0.4	0.42	105		30	150
		Nitrobenzene-d5	0.4	0.36	91		55	111
		2-Fluorobiphenyl	0.4	0.35	87		53	106
		Terphenyl-d14	0.4	0.55	137	*	58	132
Q3371-03	VPB184-HYD-20251010	2-Methylnaphthalene-d10	0.4	0.016	4	*	30	150
		Fluoranthene-d10	0.4	0.042	10	*	30	150
		Nitrobenzene-d5	0.4	0.34	86		55	111
		2-Fluorobiphenyl	0.4	0.34	84		53	106
		Terphenyl-d14	0.4	0.55	137	*	58	132
Q3371-05	BP-VPB-184-GW-920-922	2-Methylnaphthalene-d10	0.4	0.29	73		30	150
		Fluoranthene-d10	0.4	0.41	101		30	150
		Nitrobenzene-d5	0.4	0.35	88		55	111
		2-Fluorobiphenyl	0.4	0.33	83		53	106
		Terphenyl-d14	0.4	0.55	137	*	58	132
Q3371-06	VPB184-HYD-20251016	2-Methylnaphthalene-d10	0.4	0.014	4	*	30	150
		Fluoranthene-d10	0.4	0.043	11	*	30	150
		Nitrobenzene-d5	0.4	0.31	77		55	111
		2-Fluorobiphenyl	0.4	0.32	79		53	106
		Terphenyl-d14	0.4	0.52	131		58	132

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3371

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN038103.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	RPD	Limits		RPD
								Qual	Low	High	
PB170167BS	1,4-Dioxane	0.4	0.38	ug/L	95				70	130	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3371

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN038104.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	RPD	Limits		RPD
								Qual	Low	High	
PB170167BSD	1,4-Dioxane	0.4	0.35	ug/L	88	8			70	130	20

4B

SEMIVOLATILE METHOD BLANK SUMMARY

Client ID

PB170167BL

Lab Name: Alliance

Contract: TETR06

Lab Code: ACE

SDG NO.: Q3371

Lab File ID: BN038102.D

Lab Sample ID: PB170167BL

Instrument ID: BNA_N

Date Extracted: 10/20/2025

Matrix: (soil/water) Water

Date Analyzed: 10/20/2025

Level: (low/med) LOW

Time Analyzed: 17:56

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
PB170167BS	PB170167BS	BN038103.D	10/20/2025
PB170167BSD	PB170167BSD	BN038104.D	10/20/2025
BP-VPB-184-EB-20251016	Q3371-02	BN038105.D	10/20/2025
VPB184-HYD-20251010	Q3371-03	BN038106.D	10/20/2025
BP-VPB-184-GW-920-922	Q3371-05	BN038107.D	10/20/2025
VPB184-HYD-20251016	Q3371-06	BN038108.D	10/20/2025

COMMENTS: _____

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance
Lab Code: ACE
Lab File ID: BN037860.D
Instrument ID: BNA_N

Contract: TETR06
SDG NO.: Q3371
DFTPP Injection Date: 09/23/2025
DFTPP Injection Time: 11:33

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
68	Less than 2.0% of mass 69	0.7 (1.8) 1
69	Mass 69 relative abundance	100
70	Less than 2.0% of mass 69	0.2 (0.6) 1
197	Less than 2.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100
199	5.0 to 9.0% of mass 198	6.9
365	Greater than 1% of mass 198	4.2
441	Present, but less than mass 443	72.6
442	Greater than 50% of mass 198	100
443	15.0 - 24.0% of mass 442	17.7 (20.8) 2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
SSTDICC0.1	SSTDICC0.1	BN037861.D	09/23/2025	12:13
SSTDICC0.2	SSTDICC0.2	BN037862.D	09/23/2025	12:49
SSTDICCC0.4	SSTDICCC0.4	BN037863.D	09/23/2025	13:25
SSTDICC0.8	SSTDICC0.8	BN037864.D	09/23/2025	14:02
SSTDICC1.6	SSTDICC1.6	BN037865.D	09/23/2025	14:38
SSTDICC3.2	SSTDICC3.2	BN037866.D	09/23/2025	15:15
SSTDICC5.0	SSTDICC5.0	BN037867.D	09/23/2025	15:51

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance
Lab Code: ACE
Lab File ID: BN038100.D
Instrument ID: BNA_N

Contract: TETR06
SDG NO.: Q3371
DFTPP Injection Date: 10/20/2025
DFTPP Injection Time: 16:41

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
68	Less than 2.0% of mass 69	0.7 (1.8) 1
69	Mass 69 relative abundance	100
70	Less than 2.0% of mass 69	0.2 (0.5) 1
197	Less than 2.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100
199	5.0 to 9.0% of mass 198	6.8
365	Greater than 1% of mass 198	3.6
441	Present, but less than mass 443	82.8
442	Greater than 50% of mass 198	100
443	15.0 - 24.0% of mass 442	17 (19.6) 2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
SSTDCCC0.4	SSTDCCC0.4	BN038101.D	10/20/2025	17:20
PB170167BL	PB170167BL	BN038102.D	10/20/2025	17:56
PB170167BS	PB170167BS	BN038103.D	10/20/2025	18:33
PB170167BSD	PB170167BSD	BN038104.D	10/20/2025	19:09
BP-VPB-184-EB-20251016	Q3371-02	BN038105.D	10/20/2025	19:45
VPB184-HYD-20251010	Q3371-03	BN038106.D	10/20/2025	20:21
BP-VPB-184-GW-920-922	Q3371-05	BN038107.D	10/20/2025	20:58
VPB184-HYD-20251016	Q3371-06	BN038108.D	10/20/2025	21:34
SSTDCCC0.4EC	SSTDCCC0.4	BN038109.D	10/20/2025	22:10

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3371

Client ID : SSTDCCC0.4

Date Analyzed: 10/20/2025

Lab File ID: BN038101.D

Time Analyzed: 17:20

Instrument ID: BNA_N

GC Column: ZB-GR ID: 0.25 (mm)

	IS1 (DCB) AREA #	RT #	IS2 (NPT) AREA #	RT #	IS3 (ANT) AREA #	RT #
12 HOUR STD	8494	7.789	23762	10.58	12771	14.44
UPPER LIMIT	16988	8.289	47524	11.084	25542	14.941
LOWER LIMIT	4247	7.289	11881	10.084	6385.5	13.941
EPA SAMPLE NO.						
01 PB170167BL	8486	7.79	21914	10.59	10397	14.45
02 PB170167BS	7519	7.79	19915	10.58	9572	14.44
03 PB170167BSD	15910	7.79	42753	10.58	20532	14.44
04 BP-VPB-184-EB-20251016	6197	7.79	16002	10.58	8401	14.44
05 VPB184-HYD-20251010	7049	7.79	18584	10.58	9633	14.44
06 BP-VPB-184-GW-920-922	7237	7.79	19084	10.58	10109	14.44
07 VPB184-HYD-20251016	7491	7.79	19850	10.58	10193	14.44

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.

* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3371

Client ID: SSTDCCC0.4

Date Analyzed: 10/20/2025

Lab File ID: BN038101.D

Time Analyzed: 17:20

Instrument ID: BNA_N

GC Column: ZB-GR ID: 0.25 (mm)

	IS4 (PHN) AREA #	RT #	IS5 (CRY) AREA #	RT #	IS6 (PRY) AREA #	RT #
12 HOUR STD	22871	17.198	16306	21.384	14546	23.701
UPPER LIMIT	45742	17.698	32612	21.884	29092	24.201
LOWER LIMIT	11435.5	16.698	8153	20.884	7273	23.201
EPA SAMPLE NO.						
01 PB170167BL	17174	17.20	12295	21.38	12453	23.70
02 PB170167BS	16635	17.20	10986	21.38	10158	23.70
03 PB170167BSD	36706	17.20	24714	21.38	22637	23.70
04 BP-VPB-184-EB-20251016	17071	17.20	15030	21.38	14485	23.70
05 VPB184-HYD-20251010	18045	17.20	14009	21.38	12701	23.70
06 BP-VPB-184-GW-920-922	20161	17.20	17522	21.38	17404	23.69
07 VPB184-HYD-20251016	19689	17.20	16604	21.38	14840	23.69

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.

* Values outside of QC limits.



QC SAMPLE DATA

Report of Analysis

Client: Tetra Tech NUS, Inc.
Project: NWIRP Bethpage 112G08005-WE13
Client Sample ID: PB170167BL
Lab Sample ID: PB170167BL
Analytical Method: SW8270ESIM Level : LOW
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL
Prep Method : Prep Date: 10/20/25

Date Collected:
Date Received:
SDG No.: Q3371
Matrix: Water
% Solid: 0
Test: SVOC-SIMGroup1

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.20	U	1	0.070	0.20	0.20	ug/L	10/20/25 17:56	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.34			30 - 150		85%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.33			30 - 150		83%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.36			55 - 111		90%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.39			53 - 106		98%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.45			58 - 132		113%	SPK: 0.4		
INTERNAL STANDARDS										
		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	8490								
1146-65-2	Naphthalene-d8	21900								
15067-26-2	Acenaphthene-d10	10400								
1517-22-2	Phenanthrene-d10	17200								
1719-03-5	Chrysene-d12	12300								
1520-96-3	Perylene-d12	12500								

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Report of Analysis

Client: Tetra Tech NUS, Inc.
Project: NWIRP Bethpage 112G08005-WE13
Client Sample ID: PB170167BS
Lab Sample ID: PB170167BS
Analytical Method: SW8270ESIM Level : LOW
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL
Prep Method : Prep Date: 10/20/25

Date Collected:
Date Received:
SDG No.: Q3371
Matrix: Water
% Solid: 0
Test: SVOC-SIMGroup1

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.38		1	0.070	0.20	0.20	ug/L	10/20/25 18:33	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.40			30 - 150		99%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.34			30 - 150		85%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.42			55 - 111		104%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.44	*		53 - 106		111%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.51			58 - 132		127%	SPK: 0.4		
INTERNAL STANDARDS		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	7520								
1146-65-2	Naphthalene-d8	19900								
15067-26-2	Acenaphthene-d10	9570								
1517-22-2	Phenanthrene-d10	16600								
1719-03-5	Chrysene-d12	11000								
1520-96-3	Perylene-d12	10200								

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J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Report of Analysis

Client: Tetra Tech NUS, Inc.
Project: NWIRP Bethpage 112G08005-WE13
Client Sample ID: PB170167BSD
Lab Sample ID: PB170167BSD
Analytical Method: SW8270ESIM Level : LOW
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL
Prep Method : Prep Date: 10/20/25

Date Collected:
Date Received:
SDG No.: Q3371
Matrix: Water
% Solid: 0
Test: SVOC-SIMGroup1

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOD	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS										
123-91-1	1,4-Dioxane	0.35		1	0.070	0.20	0.20	ug/L	10/20/25 19:09	PB170167
SURROGATES										
7297-45-2	2-Methylnaphthalene-d10	0.36			30 - 150		90%	SPK: 0.4		
93951-69-0	Fluoranthene-d10	0.31			30 - 150		78%	SPK: 0.4		
4165-60-0	Nitrobenzene-d5	0.38			55 - 111		95%	SPK: 0.4		
321-60-8	2-Fluorobiphenyl	0.42			53 - 106		105%	SPK: 0.4		
1718-51-0	Terphenyl-d14	0.47			58 - 132		117%	SPK: 0.4		
INTERNAL STANDARDS										
		Area Count								
3855-82-1	1,4-Dichlorobenzene-d4	15900								
1146-65-2	Naphthalene-d8	42800								
15067-26-2	Acenaphthene-d10	20500								
1517-22-2	Phenanthrene-d10	36700								
1719-03-5	Chrysene-d12	24700								
1520-96-3	Perylene-d12	22600								

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products



CALIBRATION SUMMARY

Method Path : Z:\svoasrv\HPCHEM1\BNA_N\Methods\
 Method File : 8270-SIM-BN092325.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Wed Sep 24 11:00:01 2025
 Response Via : Initial Calibration

Calibration Files

0.1 =BN037861.D 0.2 =BN037862.D 0.4 =BN037863.D 0.8 =BN037864.D 1.6 =BN037865.D 3.2 =BN037866.D 5 =BN037867.D

	Compound	0.1	0.2	0.4	0.8	1.6	3.2	5	Avg	%RSD

1) I	1,4-Dichlorobenzen...	-----ISTD-----								
2)	1,4-Dioxane	0.413	0.410	0.403	0.425	0.410	0.376	0.406	4.04	
3)	n-Nitrosodimet...	0.541	0.517	0.526	0.561	0.549	0.547	0.540	3.03	
4) S	2-Fluorophenol	0.944	0.921	0.856	0.839	0.908	0.923	0.998	0.913	5.84
5) S	Phenol-d6	1.190	1.200	1.118	1.088	1.211	1.236	1.353	1.199	7.15
6)	bis(2-Chloroet...	1.133	1.096	1.082	1.066	1.150	1.124	1.143	1.113	2.89
7) I	Naphthalene-d8	-----ISTD-----								
8) S	Nitrobenzene-d5	0.299	0.289	0.289	0.290	0.310	0.320	0.338	0.305	6.14
9)	Naphthalene	1.105	1.089	1.048	1.051	1.123	1.137	1.160	1.102	3.84
10)	Hexachlorobuta...	0.198	0.184	0.182	0.183	0.192	0.188	0.188	0.188	2.94
11) SURR2	Methylnaphth...	0.515	0.518	0.499	0.504	0.553	0.587	0.714	0.556	13.74
12)	2-Methylnaphth...	0.678	0.681	0.668	0.682	0.754	0.775	0.802	0.720	7.68
13) I	Acenaphthene-d10	-----ISTD-----								
14) S	2,4,6-Tribromo...	0.148	0.142	0.135	0.139	0.163	0.184	0.152	12.15	
15) S	2-Fluorobiphenyl	1.660	1.631	1.559	1.547	1.654	1.650	1.762	1.638	4.36
16)	Acenaphthylene	1.698	1.721	1.663	1.680	1.907	1.951	2.094	1.816	9.23
17)	Acenaphthene	1.262	1.230	1.198	1.208	1.351	1.371	1.437	1.294	7.16
18)	Fluorene	1.455	1.455	1.452	1.466	1.664	1.717	1.748	1.565	8.77
19) I	Phenanthrene-d10	-----ISTD-----								
20)	4,6-Dinitro-2-...	0.041	0.044	0.046	0.058	0.071	0.084	0.057	29.75	
21)	4-Bromophenyl-...	0.252	0.247	0.244	0.244	0.272	0.275	0.291	0.261	7.07
22)	Hexachlorobenzene	0.317	0.312	0.305	0.300	0.324	0.323	0.332	0.316	3.53
23)	Atrazine	0.180	0.181	0.172	0.178	0.205	0.224	0.249	0.199	14.70
24)	Pentachlorophenol	0.101	0.092	0.097	0.116	0.136	0.108	16.50		
25)	Phenanthrene	1.301	1.271	1.230	1.224	1.358	1.379	1.452	1.316	6.37
26)	Anthracene	1.034	1.035	1.023	1.057	1.209	1.275	1.361	1.142	12.10
27) SURR	Fluoranthene-d10	0.880	0.918	0.887	0.900	1.004	1.085	1.370	1.006	17.59
28)	Fluoranthene	1.278	1.315	1.311	1.358	1.555	1.617	1.713	1.450	12.07
29) I	Chrysene-d12	-----ISTD-----								
30)	Pyrene	1.576	1.600	1.521	1.476	1.607	1.595	1.679	1.579	4.12
31) S	Terphenyl-d14	0.799	0.791	0.753	0.735	0.798	0.809	0.893	0.797	6.34
32)	Benzo(a)anthra...	1.376	1.359	1.294	1.285	1.423	1.483	1.567	1.398	7.27
33)	Chrysene	1.508	1.526	1.462	1.446	1.559	1.512	1.569	1.512	3.03
34)	Bis(2-ethylhex...	0.592	0.487	0.476	0.518	0.569	0.676	0.553	13.57	
35) I	Perylene-d12	-----ISTD-----								

Method Path : Z:\svoasrv\HPCHEM1\BNA_N\Methods\

Method File : 8270-SIM-BN092325.M

36)	Indeno(1,2,3-c...	1.715	1.672	1.658	1.707	1.887	1.996	2.159	1.828	10.52
37)	Benzo(b)fluora...	1.599	1.606	1.641	1.590	1.810	1.848	1.981	1.725	8.95
38)	Benzo(k)fluora...	1.620	1.586	1.597	1.628	1.875	1.917	1.926	1.736	9.26
39) C	Benzo(a)pyrene	1.254	1.296	1.231	1.254	1.413	1.492	1.610	1.364	10.61
40)	Dibenzo(a,h)an...	1.285	1.283	1.303	1.333	1.495	1.587	1.705	1.427	11.87
41)	Benzo(g,h,i)pe...	1.407	1.381	1.404	1.397	1.521	1.613	1.773	1.499	9.85

(#) = Out of Range

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06
Lab Code: ACE SDG No.: Q3371
Instrument ID: BNA_N Calibration Date/Time: 10/20/2025 17:20
Lab File ID: BN038101.D Init. Calib. Date(s): 09/23/2025 09/23/2025
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 12:13 15:51
GC Column: ZB-GR ID: 0.25 (mm)

COMPOUND	RRF	RRF0.4	MIN RRF	%D	MAX%D
2-Methylnaphthalene-d10	0.556	0.539		-3.1	20.0
Fluoranthene-d10	1.006	0.870		-13.5	20.0
2-Fluorophenol	0.913	0.879		-3.7	20.0
Phenol-d6	1.199	1.019		-15.0	20.0
Nitrobenzene-d5	0.305	0.277		-9.2	20.0
2-Fluorobiphenyl	1.638	1.602		-2.2	20.0
2,4,6-Tribromophenol	0.152	0.135		-11.2	20.0
Terphenyl-d14	0.797	0.829		4.0	20.0
1,4-Dioxane	0.406	0.472		16.3	20.0

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>TETR06</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3371</u>
Instrument ID:	<u>BNA_N</u>	Calibration Date/Time:	<u>10/20/2025 22:10</u>
Lab File ID:	<u>BN038109.D</u>	Init. Calib. Date(s):	<u>09/23/2025 09/23/2025</u>
EPA Sample No.:	<u>SSTDCCC0.4EC</u>	Init. Calib. Time(s):	<u>12:13 15:51</u>
GC Column:	<u>ZB-GR</u>	ID:	<u>0.25</u> (mm)

COMPOUND	RRF	RRF0.4	MIN RRF	%D	MAX%D
2-Methylnaphthalene-d10	0.556	0.519		-6.7	50.0
Fluoranthene-d10	1.006	0.890		-11.5	50.0
2-Fluorophenol	0.913	0.912		-0.1	50.0
Phenol-d6	1.199	1.015		-15.3	50.0
Nitrobenzene-d5	0.305	0.273		-10.5	50.0
2-Fluorobiphenyl	1.638	1.528		-6.7	50.0
2,4,6-Tribromophenol	0.152	0.126		-17.1	50.0
Terphenyl-d14	0.797	0.855		7.3	50.0
1,4-Dioxane	0.406	0.486		19.7	50.0

All other compounds must meet a minimum RRF of 0.010.