

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: _____

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG No.: BN110724

Instrument ID: BNA_N Calibration Date/Time: 11/7/2024 1:16:58

Lab File ID: BN034896.D Init. Calib. Date(s): _____

EPA Sample No.: SSTDCCC0.4EC Init. Calib. Time(s): _____

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

COMPOUND	RRF	RRF0.4	MIN RRF	%D	MAX%D
2-Methylnaphthalene-d10	0.545	0.495		-9.174	50.0
Fluoranthene-d10	0.902	0.847		-6.098	50.0
n-Nitrosodimethylamine	0.682	0.642		-5.865	50.0
2-Fluorophenol	1.115	1.012		-9.238	50.0
Phenol-d6	1.480	1.302		-12.027	50.0
bis(2-Chloroethyl)ether	1.277	1.170		-8.379	50.0
Nitrobenzene-d5	0.312	0.278		-10.897	50.0
Naphthalene	1.110	1.034		-6.847	50.0
Hexachlorobutadiene	0.177	0.167		-5.65	50.0
2-Methylnaphthalene	0.679	0.615		-9.426	50.0
2-Fluorobiphenyl	1.690	1.498		-11.361	50.0
Acenaphthylene	1.929	1.660		-13.945	50.0
Acenaphthene	1.335	1.165		-12.734	50.0
Fluorene	1.662	1.489		-10.409	50.0
4,6-Dinitro-2-methylphenol	0.044	0.031		-29.545	50.0
2,4,6-Tribromophenol	0.118	0.092		-22.034	50.0
4-Bromophenyl-phenylether	0.213	0.194		-8.92	50.0
Hexachlorobenzene	0.257	0.246		-4.28	50.0
Atrazine	0.154	0.135		-12.338	50.0
Pentachlorophenol	0.077	0.057		-25.974	50.0
Phenanthrene	1.227	1.155		-5.868	50.0
Anthracene	1.058	0.971		-8.223	50.0
Fluoranthene	1.291	1.226		-5.035	50.0
Pyrene	2.025	1.965		-2.963	50.0
Terphenyl-d14	0.749	0.727		-2.937	50.0
Benzo(a)anthracene	1.559	1.484		-4.811	50.0
Chrysene	1.650	1.633		-1.03	50.0
Bis(2-ethylhexyl)phthalate	0.895	0.718		-19.777	50.0
Benzo(b)fluoranthene	1.757	1.779		1.252	50.0
Benzo(k)fluoranthene	1.828	1.811		-0.93	50.0
Benzo(a)pyrene	1.396	1.311		-6.089	50.0
Indeno(1,2,3-cd)pyrene	1.782	1.585		-11.055	50.0
Dibenzo(a,h)anthracene	1.379	1.203		-12.763	50.0
Benzo(g,h,i)perylene	1.464	1.349		-7.855	50.0
1,4-Dioxane	0.505	0.471		-6.733	50.0

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: _____

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG No.: BN110724

Instrument ID: BNA_N Calibration Date/Time: 11/7/2024 11:24

Lab File ID: BN034887.D Init. Calib. Date(s): _____

EPA Sample No.: SSTDICCC0.4 Init. Calib. Time(s): _____

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

COMPOUND	RRF	RRF0.4	MIN RRF	%D	MAX%D
2-Methylnaphthalene-d10	0.545	0.491		-9.908	20.0
Fluoranthene-d10	0.902	0.801		-11.197	20.0
n-Nitrosodimethylamine	0.682	0.621		-8.944	20.0
2-Fluorophenol	1.115	1.007		-9.686	20.0
Phenol-d6	1.480	1.299		-12.23	20.0
bis(2-Chloroethyl) ether	1.277	1.170		-8.379	20.0
Nitrobenzene-d5	0.312	0.280		-10.256	20.0
Naphthalene	1.110	1.028		-7.387	20.0
Hexachlorobutadiene	0.177	0.167		-5.65	20.0
2-Methylnaphthalene	0.679	0.612		-9.867	20.0
2-Fluorobiphenyl	1.690	1.540		-8.876	20.0
Acenaphthylene	1.929	1.672		-13.323	20.0
Acenaphthene	1.335	1.176		-11.91	20.0
Fluorene	1.662	1.463		-11.974	20.0
4,6-Dinitro-2-methylphenol	0.044	0.032		-27.273	20.0
2,4,6-Tribromophenol	0.118	0.087		-26.271	20.0
4-Bromophenyl-phenylether	0.213	0.202		-5.164	20.0
Hexachlorobenzene	0.257	0.252		-1.946	20.0
Atrazine	0.154	0.133		-13.636	20.0
Pentachlorophenol	0.077	0.055		-28.571	20.0
Phenanthrene	1.227	1.162		-5.297	20.0
Anthracene	1.058	0.963		-8.979	20.0
Fluoranthene	1.291	1.135		-12.084	20.0
Pyrene	2.025	1.990		-1.728	20.0
Terphenyl-d14	0.749	0.743		-0.801	20.0
Benzo(a)anthracene	1.559	1.423		-8.724	20.0
Chrysene	1.650	1.583		-4.061	20.0
Bis(2-ethylhexyl)phthalate	0.895	0.720		-19.553	20.0
Benzo(b)fluoranthene	1.757	1.752		-0.285	20.0
Benzo(k)fluoranthene	1.828	1.724		-5.689	20.0
Benzo(a)pyrene	1.396	1.339		-4.083	20.0
Indeno(1,2,3-cd)pyrene	1.782	1.680		-5.724	20.0
Dibenzo(a,h)anthracene	1.379	1.304		-5.439	20.0
Benzo(g,h,i)perylene	1.464	1.483		1.298	20.0
1,4-Dioxane	0.505	0.474		-6.139	20.0

All other compounds must meet a minimum RRF of 0.010.

Report of Analysis

Client:		Date Collected:	
Project:		Date Received:	
Client Sample ID:	SSTDCCC0.4	SDG No.:	BN110724
Lab Sample ID:	SSTDCCC0.4	Matrix:	Water
Analytical Method:	8270-Modified	% Moisture:	100
Sample Wt/Vol:	1000 Units: mL	Final Vol:	1000000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :		Decanted :	
Injection Volume :		Level :	LOW
	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034884.D	1		11/7/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	370		30	200	200	ug/L
111-44-4	bis(2-Chloroethyl)ether	370		28	100	100	ug/L
91-20-3	Naphthalene	370		24	100	100	ug/L
87-68-3	Hexachlorobutadiene	380		26	100	100	ug/L
91-57-6	2-Methylnaphthalene	360		30	100	100	ug/L
208-96-8	Acenaphthylene	350		21	100	100	ug/L
83-32-9	Acenaphthene	350		20	100	100	ug/L
Total PAH	Total PAH	6010		456	100	1600	ug/L
86-73-7	Fluorene	350		18	100	100	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	330		140	200	200	ug/L
101-55-3	4-Bromophenyl-phenylether	380		26	100	100	ug/L
118-74-1	Hexachlorobenzene	400		29	100	100	ug/L
1912-24-9	Atrazine	330		23	100	100	ug/L
87-86-5	Pentachlorophenol	330		90	200	200	ug/L
85-01-8	Phenanthrene	370		20	100	100	ug/L
120-12-7	Anthracene	360		24	100	100	ug/L
206-44-0	Fluoranthene	340		23	100	100	ug/L
129-00-0	Pyrene	410		22	100	100	ug/L
56-55-3	Benzo(a)anthracene	360		22	100	100	ug/L
218-01-9	Chrysene	390		26	100	100	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	320		140	200	200	ug/L
205-99-2	Benzo(b)fluoranthene	410		32	100	100	ug/L
207-08-9	Benzo(k)fluoranthene	380		34	100	100	ug/L
50-32-8	Benzo(a)pyrene	380		56	100	100	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	380		38	100	100	ug/L
53-70-3	Dibenzo(a,h)anthracene	390		39	100	100	ug/L
191-24-2	Benzo(g,h,i)perylene	420		37	100	100	ug/L
123-91-1	1,4-Dioxane	370		68	200	200	ug/L

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	SSTDCCC0.4			SDG No.:	BN110724		
Lab Sample ID:	SSTDCCC0.4			Matrix:	Water		
Analytical Method:	8270-Modified			% Moisture:	100		
Sample Wt/Vol:	1000	Units:	mL	Final Vol:	1000000	uL	
Soil Aliquot Vol:			uL	Test:			
Extraction Type :		Decanted :		Level :	LOW		
Injection Volume :		GPC Factor :		GPC Cleanup :		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034884.D	1		11/7/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	362	*	20-139		90500	SPK: -0.4
93951-69-0	Fluoranthene-d10	344	*	30-150		86000	SPK: -0.4
367-12-4	2-Fluorophenol	355	*	10-100		88750	SPK: -0.4
13127-88-3	Phenol-d6	348	*	10-100		87000	SPK: -0.4
4165-60-0	Nitrobenzene-d5	369	*	27-123		92250	SPK: -0.4
321-60-8	2-Fluorobiphenyl	367	*	34-132		91750	SPK: -0.4
118-79-6	2,4,6-Tribromophenol	326	*	10-131		81500	SPK: -0.4
1718-51-0	Terphenyl-d14	405	*	35-157		101250	SPK: -0.4
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	6014	7.575				
1146-65-2	Naphthalene-d8	17441	10.34				
15067-26-2	Acenaphthene-d10	8024	14.208				
1517-22-2	Phenanthrene-d10	15173	16.952				
1719-03-5	Chrysene-d12	8034	21.149				
1520-96-3	Perylene-d12	6346	23.315				

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBN110724			SDG No.:	BN110724		
Lab Sample ID:	SSTDICV0.4			Matrix:	Water		
Analytical Method:	8270-Modified			% Moisture:	100		
Sample Wt/Vol:	1000	Units:	mL	Final Vol:	1000000	uL	
Soil Aliquot Vol:			uL	Test:			
Extraction Type :		Decanted :		Level :	LOW		
Injection Volume :		GPC Factor :		GPC Cleanup :		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034892.D	1		11/7/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	340		30	200	200	ug/L
111-44-4	bis(2-Chloroethyl)ether	360		28	100	100	ug/L
91-20-3	Naphthalene	380		24	100	100	ug/L
87-68-3	Hexachlorobutadiene	380		26	100	100	ug/L
91-57-6	2-Methylnaphthalene	380		30	100	100	ug/L
208-96-8	Acenaphthylene	350		21	100	100	ug/L
83-32-9	Acenaphthene	360		20	100	100	ug/L
Total PAH	Total PAH	5810		456	100	1600	ug/L
86-73-7	Fluorene	360		18	100	100	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	350		140	200	200	ug/L
101-55-3	4-Bromophenyl-phenylether	380		26	100	100	ug/L
118-74-1	Hexachlorobenzene	390		29	100	100	ug/L
1912-24-9	Atrazine	360		23	100	100	ug/L
87-86-5	Pentachlorophenol	370		90	200	200	ug/L
85-01-8	Phenanthrene	380		20	100	100	ug/L
120-12-7	Anthracene	370		24	100	100	ug/L
206-44-0	Fluoranthene	380		23	100	100	ug/L
129-00-0	Pyrene	370		22	100	100	ug/L
56-55-3	Benzo(a)anthracene	370		22	100	100	ug/L
218-01-9	Chrysene	390		26	100	100	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	330		140	200	200	ug/L
205-99-2	Benzo(b)fluoranthene	380		32	100	100	ug/L
207-08-9	Benzo(k)fluoranthene	360		34	100	100	ug/L
50-32-8	Benzo(a)pyrene	350		56	100	100	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	330		38	100	100	ug/L
53-70-3	Dibenzo(a,h)anthracene	330		39	100	100	ug/L
191-24-2	Benzo(g,h,i)perylene	350		37	100	100	ug/L
123-91-1	1,4-Dioxane	350		68	200	200	ug/L

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBN110724			SDG No.:	BN110724		
Lab Sample ID:	SSTDICV0.4			Matrix:	Water		
Analytical Method:	8270-Modified			% Moisture:	100		
Sample Wt/Vol:	1000	Units:	mL	Final Vol:	1000000	uL	
Soil Aliquot Vol:			uL	Test:			
Extraction Type :		Decanted :		Level :	LOW		
Injection Volume :		GPC Factor :		GPC Cleanup :		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034892.D	1		11/7/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	379	*	20-139		94750	SPK: -0.4
93951-69-0	Fluoranthene-d10	383	*	30-150		95750	SPK: -0.4
367-12-4	2-Fluorophenol	354	*	10-100		88500	SPK: -0.4
13127-88-3	Phenol-d6	348	*	10-100		87000	SPK: -0.4
4165-60-0	Nitrobenzene-d5	345	*	27-123		86250	SPK: -0.4
321-60-8	2-Fluorobiphenyl	359	*	34-132		89750	SPK: -0.4
118-79-6	2,4,6-Tribromophenol	383	*	10-131		95750	SPK: -0.4
1718-51-0	Terphenyl-d14	385	*	35-157		96250	SPK: -0.4
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	5596	7.575				
1146-65-2	Naphthalene-d8	16947	10.34				
15067-26-2	Acenaphthene-d10	8617	14.208				
1517-22-2	Phenanthrene-d10	17630	16.952				
1719-03-5	Chrysene-d12	11730	21.149				
1520-96-3	Perylene-d12	10292	23.315				

Report of Analysis

Client:		Date Collected:	10/17/2024 12:00:00 AM
Project:		Date Received:	10/17/2024 12:00:00 AM
Client Sample ID:	PB164229BL	SDG No.:	BN110724
Lab Sample ID:	PB164229BL	Matrix:	Water
Analytical Method:	8270-Modified	% Moisture:	100
Sample Wt/Vol:	1000 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034893.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164229

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	0.03	U	0.03	0.2	0.2	ug/L
111-44-4	bis(2-Chloroethyl)ether	0.03	U	0.03	0.1	0.1	ug/L
91-20-3	Naphthalene	0.02	U	0.02	0.1	0.1	ug/L
87-68-3	Hexachlorobutadiene	0.03	U	0.03	0.1	0.1	ug/L
91-57-6	2-Methylnaphthalene	0.03	U	0.03	0.1	0.1	ug/L
208-96-8	Acenaphthylene	0.02	U	0.02	0.1	0.1	ug/L
83-32-9	Acenaphthene	0.02	U	0.02	0.1	0.1	ug/L
Total PAH	Total PAH	0.45	U	0.45	0.1	1.6	ug/L
86-73-7	Fluorene	0.02	U	0.02	0.1	0.1	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	0.14	U	0.14	0.2	0.2	ug/L
101-55-3	4-Bromophenyl-phenylether	0.03	U	0.03	0.1	0.1	ug/L
118-74-1	Hexachlorobenzene	0.03	U	0.03	0.1	0.1	ug/L
1912-24-9	Atrazine	0.02	U	0.02	0.1	0.1	ug/L
87-86-5	Pentachlorophenol	0.09	U	0.09	0.2	0.2	ug/L
85-01-8	Phenanthrene	0.02	U	0.02	0.1	0.1	ug/L
120-12-7	Anthracene	0.02	U	0.02	0.1	0.1	ug/L
206-44-0	Fluoranthene	0.02	U	0.02	0.1	0.1	ug/L
129-00-0	Pyrene	0.02	U	0.02	0.1	0.1	ug/L
56-55-3	Benzo(a)anthracene	0.02	U	0.02	0.1	0.1	ug/L
218-01-9	Chrysene	0.03	U	0.03	0.1	0.1	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	0.14	U	0.14	0.2	0.2	ug/L
205-99-2	Benzo(b)fluoranthene	0.03	U	0.03	0.1	0.1	ug/L
207-08-9	Benzo(k)fluoranthene	0.03	U	0.03	0.1	0.1	ug/L
50-32-8	Benzo(a)pyrene	0.06	U	0.06	0.1	0.1	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	0.04	U	0.04	0.1	0.1	ug/L
53-70-3	Dibenzo(a,h)anthracene	0.04	U	0.04	0.1	0.1	ug/L
191-24-2	Benzo(g,h,i)perylene	0.04	U	0.04	0.1	0.1	ug/L
123-91-1	1,4-Dioxane	0.07	U	0.07	0.2	0.2	ug/L

Report of Analysis

Client:		Date Collected:	10/17/2024 12:00:00 AM
Project:		Date Received:	10/17/2024 12:00:00 AM
Client Sample ID:	PB164229BL	SDG No.:	BN110724
Lab Sample ID:	PB164229BL	Matrix:	Water
Analytical Method:	8270-Modified	% Moisture:	100
Sample Wt/Vol:	1000 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034893.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164229

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	0.379		20-139		95	SPK: -0.4
93951-69-0	Fluoranthene-d10	0.401		30-150		100	SPK: -0.4
367-12-4	2-Fluorophenol	0.364		10-100		91	SPK: -0.4
13127-88-3	Phenol-d6	0.358		10-100		90	SPK: -0.4
4165-60-0	Nitrobenzene-d5	0.38		27-123		95	SPK: -0.4
321-60-8	2-Fluorobiphenyl	0.405		34-132		101	SPK: -0.4
118-79-6	2,4,6-Tribromophenol	0.282		10-131		70	SPK: -0.4
1718-51-0	Terphenyl-d14	0.519		35-157		130	SPK: -0.4
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	6231	7.575				
1146-65-2	Naphthalene-d8	17894	10.34				
15067-26-2	Acenaphthene-d10	7932	14.208				
1517-22-2	Phenanthrene-d10	16492	16.952				
1719-03-5	Chrysene-d12	9359	21.149				
1520-96-3	Perylene-d12	7338	23.318				

Report of Analysis

Client:		Date Collected:	10/9/2024 12:00:00 AM
Project:		Date Received:	10/9/2024 12:00:00 AM
Client Sample ID:	LOQ-SOIL-02-QT4-2024	SDG No.:	BN110724
Lab Sample ID:	P4368-02	Matrix:	Solid
Analytical Method:	8270-Modified	% Moisture:	0
Sample Wt/Vol:	30.01 Units: g	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034894.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164230

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	6.8		1.3	6.6	6.6	ug/Kg
111-44-4	bis(2-Chloroethyl)ether	7.3		0.85	3.3	3.3	ug/Kg
91-20-3	Naphthalene	7.2		0.69	3.3	3.3	ug/Kg
87-68-3	Hexachlorobutadiene	7		0.69	3.3	3.3	ug/Kg
91-57-6	2-Methylnaphthalene	7.1		0.8	3.3	3.3	ug/Kg
208-96-8	Acenaphthylene	7		0.63	3.3	3.3	ug/Kg
83-32-9	Acenaphthene	6.8		0.55	3.3	3.3	ug/Kg
86-73-7	Fluorene	6.8		0.6	3.3	3.3	ug/Kg
534-52-1	4,6-Dinitro-2-methylphenol	6.1	J	4.2	6.6	6.6	ug/Kg
101-55-3	4-Bromophenyl-phenylether	6.9		0.79	3.3	3.3	ug/Kg
118-74-1	Hexachlorobenzene	7.1		0.85	3.3	3.3	ug/Kg
1912-24-9	Atrazine	6.7		0.76	3.3	3.3	ug/Kg
87-86-5	Pentachlorophenol	5.9	J	2.8	6.6	6.6	ug/Kg
85-01-8	Phenanthrene	7.3		0.63	3.3	3.3	ug/Kg
120-12-7	Anthracene	7.4		0.79	3.3	3.3	ug/Kg
206-44-0	Fluoranthene	7.1		0.67	3.3	3.3	ug/Kg
129-00-0	Pyrene	7.1		0.95	3.3	3.3	ug/Kg
56-55-3	Benzo(a)anthracene	7.3		0.76	3.3	3.3	ug/Kg
218-01-9	Chrysene	7.6		0.86	3.3	3.3	ug/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	6.4	J	1.5	6.6	6.6	ug/Kg
205-99-2	Benzo(b)fluoranthene	8		1.2	3.3	3.3	ug/Kg
207-08-9	Benzo(k)fluoranthene	7.3		0.98	3.3	3.3	ug/Kg
50-32-8	Benzo(a)pyrene	7.1		1.7	3.3	3.3	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	6.5		1.3	3.3	3.3	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	6.6		1.2	3.3	3.3	ug/Kg
191-24-2	Benzo(g,h,i)perylene	7.1		1.1	3.3	3.3	ug/Kg
123-91-1	1,4-Dioxane	7		2.2	6.6	6.6	ug/Kg

SURROGATES

Report of Analysis

Client:		Date Collected:	10/9/2024 12:00:00 AM
Project:		Date Received:	10/9/2024 12:00:00 AM
Client Sample ID:	LOQ-SOIL-02-QT4-2024	SDG No.:	BN110724
Lab Sample ID:	P4368-02	Matrix:	Solid
Analytical Method:	8270-Modified	% Moisture:	0
Sample Wt/Vol:	30.01 Units: g	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034894.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164230

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
7297-45-2	2-Methylnaphthalene-d10	0.563	*	17-161			SPK: -0
93951-69-0	Fluoranthene-d10	0.606	*	23-138			SPK: -0
367-12-4	2-Fluorophenol	0.312	*	28-122			SPK: -0
13127-88-3	Phenol-d6	0.317	*	25-125			SPK: -0
4165-60-0	Nitrobenzene-d5	0.354	*	33-121			SPK: -0
321-60-8	2-Fluorobiphenyl	0.349	*	32-121			SPK: -0
118-79-6	2,4,6-Tribromophenol	0.307	*	19-110			SPK: -0
1718-51-0	Terphenyl-d14	0.381	*	21-130			SPK: -0
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	5405	7.575				
1146-65-2	Naphthalene-d8	15883	10.34				
15067-26-2	Acenaphthene-d10	7600	14.211				
1517-22-2	Phenanthrene-d10	15509	16.957				
1719-03-5	Chrysene-d12	9605	21.142				
1520-96-3	Perylene-d12	7636	23.314				

Report of Analysis

Client:		Date Collected:	10/9/2024 12:00:00 AM
Project:		Date Received:	10/9/2024 12:00:00 AM
Client Sample ID:	LOQ-WATER-02-QT4-2024	SDG No.:	BN110724
Lab Sample ID:	P4368-08	Matrix:	Water
Analytical Method:	8270-Modified	% Moisture:	100
Sample Wt/Vol:	1000 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034895.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164229

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	0.22		0.03	0.2	0.2	ug/L
111-44-4	bis(2-Chloroethyl)ether	0.22		0.03	0.1	0.1	ug/L
91-20-3	Naphthalene	0.22		0.02	0.1	0.1	ug/L
87-68-3	Hexachlorobutadiene	0.21		0.03	0.1	0.1	ug/L
91-57-6	2-Methylnaphthalene	0.22		0.03	0.1	0.1	ug/L
208-96-8	Acenaphthylene	0.21		0.02	0.1	0.1	ug/L
83-32-9	Acenaphthene	0.2		0.02	0.1	0.1	ug/L
86-73-7	Fluorene	0.21		0.02	0.1	0.1	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	0.19	J	0.14	0.2	0.2	ug/L
101-55-3	4-Bromophenyl-phenylether	0.2		0.03	0.1	0.1	ug/L
118-74-1	Hexachlorobenzene	0.22		0.03	0.1	0.1	ug/L
1912-24-9	Atrazine	0.2		0.02	0.1	0.1	ug/L
87-86-5	Pentachlorophenol	0.18	J	0.09	0.2	0.2	ug/L
85-01-8	Phenanthrene	0.22		0.02	0.1	0.1	ug/L
120-12-7	Anthracene	0.22		0.02	0.1	0.1	ug/L
206-44-0	Fluoranthene	0.21		0.02	0.1	0.1	ug/L
129-00-0	Pyrene	0.21		0.02	0.1	0.1	ug/L
56-55-3	Benzo(a)anthracene	0.22		0.02	0.1	0.1	ug/L
218-01-9	Chrysene	0.22		0.03	0.1	0.1	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	0.19	J	0.14	0.2	0.2	ug/L
205-99-2	Benzo(b)fluoranthene	0.23		0.03	0.1	0.1	ug/L
207-08-9	Benzo(k)fluoranthene	0.22		0.03	0.1	0.1	ug/L
50-32-8	Benzo(a)pyrene	0.22		0.06	0.1	0.1	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	0.2		0.04	0.1	0.1	ug/L
53-70-3	Dibenzo(a,h)anthracene	0.2		0.04	0.1	0.1	ug/L
191-24-2	Benzo(g,h,i)perylene	0.21		0.04	0.1	0.1	ug/L
123-91-1	1,4-Dioxane	0.23		0.07	0.2	0.2	ug/L

SURROGATES

Report of Analysis

Client:		Date Collected:	10/9/2024 12:00:00 AM
Project:		Date Received:	10/9/2024 12:00:00 AM
Client Sample ID:	LOQ-WATER-02-QT4-2024	SDG No.:	BN110724
Lab Sample ID:	P4368-08	Matrix:	Water
Analytical Method:	8270-Modified	% Moisture:	100
Sample Wt/Vol:	1000 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	
Extraction Type :	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN034895.D	1	10/17/2024 12:00:00 AM	11/7/2024 12:00:00 AM	PB164229

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
7297-45-2	2-Methylnaphthalene-d10	0.568	*	20-139			SPK: -0
93951-69-0	Fluoranthene-d10	0.609	*	30-150			SPK: -0
367-12-4	2-Fluorophenol	0.317	*	10-100			SPK: -0
13127-88-3	Phenol-d6	0.32	*	10-100			SPK: -0
4165-60-0	Nitrobenzene-d5	0.358	*	27-123			SPK: -0
321-60-8	2-Fluorobiphenyl	0.35	*	34-132			SPK: -0
118-79-6	2,4,6-Tribromophenol	0.305	*	10-131			SPK: -0
1718-51-0	Terphenyl-d14	0.378	*	35-157			SPK: -0
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	4901	7.575				
1146-65-2	Naphthalene-d8	14321	10.34				
15067-26-2	Acenaphthene-d10	6899	14.208				
1517-22-2	Phenanthrene-d10	14131	16.952				
1719-03-5	Chrysene-d12	8914	21.14				
1520-96-3	Perylene-d12	7193	23.315				

Hit Summary Sheet SW-846

SDG No.: BN110724

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
Client ID : SSTDCCC0.4								
SSTDCCC0.4	SSTDCCC0.4	Water	1,4-Dioxane	370.000	68		200	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	2-Methylnaphthalene	360.000	30		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	4,6-Dinitro-2-methylphenol	330.000	140		200	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	4-Bromophenyl-phenylether	380.000	26		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Acenaphthene	350.000	20		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Acenaphthylene	350.000	21		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Anthracene	360.000	24		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Atrazine	330.000	23		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Benzo(a)anthracene	360.000	22		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Benzo(a)pyrene	380.000	56		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Benzo(b)fluoranthene	410.000	32		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Benzo(g,h,i)perylene	420.000	37		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Benzo(k)fluoranthene	380.000	34		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	bis(2-Chloroethyl)ether	370.000	28		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Bis(2-ethylhexyl)phthalate	320.000	140		200	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Chrysene	390.000	26		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Dibenzo(a,h)anthracene	390.000	39		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Fluoranthene	340.000	23		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Fluorene	350.000	18		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Hexachlorobenzene	400.000	29		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Hexachlorobutadiene	380.000	26		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Indeno(1,2,3-cd)pyrene	380.000	38		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	n-Nitrosodimethylamine	370.000	30		200	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Naphthalene	370.000	24		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Pentachlorophenol	330.000	90		200	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Phenanthrene	370.000	20		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Pyrene	410.000	22		100	ug/L
SSTDCCC0.4	SSTDCCC0.4	Water	Total PAH	6,010.000	456		1600	ug/L

Total Svoc : 15,960.00

Total Concentration: 15,960.00

Client ID : ICVBN110724

SSTDICV0.4	ICVBN110724	Water	1,4-Dioxane	350.000	68		200	ug/L
SSTDICV0.4	ICVBN110724	Water	2-Methylnaphthalene	380.000	30		100	ug/L
SSTDICV0.4	ICVBN110724	Water	4,6-Dinitro-2-methylphenol	350.000	140		200	ug/L
SSTDICV0.4	ICVBN110724	Water	4-Bromophenyl-phenylether	380.000	26		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Acenaphthene	360.000	20		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Acenaphthylene	350.000	21		100	ug/L

Hit Summary Sheet SW-846

SDG No.: BN110724

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
SSTDICV0.4	ICVBN110724	Water	Anthracene	370.000	24		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Atrazine	360.000	23		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Benzo(a)anthracene	370.000	22		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Benzo(a)pyrene	350.000	56		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Benzo(b)fluoranthene	380.000	32		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Benzo(g,h,i)perylene	350.000	37		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Benzo(k)fluoranthene	360.000	34		100	ug/L
SSTDICV0.4	ICVBN110724	Water	bis(2-Chloroethyl)ether	360.000	28		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Bis(2-ethylhexyl)phthalate	330.000	140		200	ug/L
SSTDICV0.4	ICVBN110724	Water	Chrysene	390.000	26		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Dibenzo(a,h)anthracene	330.000	39		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Fluoranthene	380.000	23		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Fluorene	360.000	18		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Hexachlorobenzene	390.000	29		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Hexachlorobutadiene	380.000	26		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Indeno(1,2,3-cd)pyrene	330.000	38		100	ug/L
SSTDICV0.4	ICVBN110724	Water	n-Nitrosodimethylamine	340.000	30		200	ug/L
SSTDICV0.4	ICVBN110724	Water	Naphthalene	380.000	24		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Pentachlorophenol	370.000	90		200	ug/L
SSTDICV0.4	ICVBN110724	Water	Phenanthrene	380.000	20		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Pyrene	370.000	22		100	ug/L
SSTDICV0.4	ICVBN110724	Water	Total PAH	5,810.000	456		1600	ug/L
Total Svoc :				15,610.00				
Total Concentration:				15,610.00				

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BN110724SAS No.: BN110724SDG No.: BN110724

Instrument ID: _____

Calibration Date(s): 11/7/2024 :Calibration Time(s): 10:02

LAB FILE ID:		RRF0.1 = BN034885.D		RRF0.2 = BN034886.D		RRF0.4 = BN034887.D		
		RRF0.8 = BN034888.D		RRF1.6 = BN034889.D		RRF3.2 = BN034890.D		
COMPOUND	RRF0.1	RRF0.2	RRF0.4	RRF0.8	RRF1.6	RRF3.2	RRF	% RSD
2-Methylnaphthalene-d10	0.527	0.519	0.491	0.579	0.577	0.548	0.545	6.3
Fluoranthene-d10	0.801	0.845	0.801	0.971	0.976	0.947	0.902	9.2
n-Nitrosodimethylamine	0.697	0.715	0.621	0.753	0.712	0.643	0.682	7.3
2-Fluorophenol	1.137	1.131	1.007	1.180	1.146	1.082	1.115	5.0
Phenol-d6	1.438	1.439	1.299	1.582	1.550	1.495	1.480	6.6
bis(2-Chloroethyl)ether	1.316	1.267	1.170	1.389	1.336	1.230	1.277	5.8
Nitrobenzene-d5	0.314	0.301	0.280	0.327	0.324	0.313	0.312	5.4
Naphthalene	1.110	1.086	1.028	1.181	1.166	1.089	1.110	4.6
Hexachlorobutadiene	0.183	0.180	0.167	0.189	0.182	0.168	0.177	4.9
2-Methylnaphthalene	0.647	0.641	0.612	0.725	0.723	0.691	0.679	6.8
2-Fluorobiphenyl	1.753	1.737	1.540	1.788	1.747	1.620	1.690	5.3
Acenaphthylene	1.833	1.826	1.672	2.034	2.040	2.003	1.929	8.0
Acenaphthene	1.292	1.281	1.176	1.430	1.412	1.352	1.335	6.8
Fluorene	1.614	1.600	1.463	1.774	1.779	1.680	1.662	6.8
4,6-Dinitro-2-methylphenol		0.033	0.032	0.040	0.044	0.052	0.044	26.2
2,4,6-Tribromophenol	0.083	0.094	0.087	0.121	0.129	0.142	0.118	26.5
4-Bromophenyl-phenylether	0.212	0.201	0.202	0.220	0.219	0.216	0.213	4.1
Hexachlorobenzene	0.262	0.255	0.252	0.268	0.260	0.252	0.257	2.6
Atrazine	0.128	0.136	0.133	0.162	0.169	0.173	0.154	13.8
Pentachlorophenol		0.055	0.055	0.074	0.079	0.092	0.077	26.6
Phenanthrene	1.218	1.168	1.162	1.303	1.276	1.239	1.227	4.2
Anthracene	0.965	0.963	0.963	1.121	1.116	1.119	1.058	8.4
Fluoranthene	1.112	1.183	1.135	1.414	1.418	1.385	1.291	10.9
Pyrene	2.180	2.005	1.990	2.107	2.033	1.945	2.025	4.5
Terphenyl-d14	0.788	0.750	0.743	0.777	0.750	0.715	0.749	3.5
Benzo (a) anthracene	1.450	1.472	1.423	1.663	1.663	1.622	1.559	6.8
Chrysene	1.632	1.632	1.583	1.768	1.709	1.625	1.650	3.9
Bis(2-ethylhexyl)phthalate	0.984	0.873	0.720	0.881	0.838	0.922	0.895	11.8
Benzo (b) fluoranthene	1.685	1.601	1.752	1.900	1.814	1.765	1.757	5.4
Benzo (k) fluoranthene	1.763	1.746	1.724	1.982	1.931	1.812	1.828	5.3
Benzo (a) pyrene	1.308	1.249	1.339	1.475	1.469	1.438	1.396	6.9
Indeno (1,2,3-cd) pyrene	1.810	1.753	1.680	1.845	1.868	1.733	1.782	3.7
Dibenzo (a,h) anthracene	1.389	1.356	1.304	1.423	1.447	1.347	1.379	3.5
Benzo (g,h,i) perylene	1.482	1.434	1.483	1.485	1.509	1.407	1.464	2.4
1,4-Dioxane	0.563	0.562	0.474	0.530	0.498	0.470	0.505	9.4

All other compounds must meet a minimum RRF of 0.010.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG NO.: BN110724

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Nov 7 2024 12:00AM

Lab File ID: BN034887.D Time Analyzed: 09:18

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	5878	7.575	16847	10.34	7757	14.21
UPPER LIMIT	11756	8.075	33694	10.84	15514	14.711
LOWER LIMIT	2939	7.075	8423.5	9.84	3878.5	13.711
EPA SAMPLE NO.						
01 SSTDCCC0.4	6014	7.58	17441	10.34	8024	14.21
02 ICVBN110724	5596	7.58	16947	10.34	8617	14.21
03 PB164229BL	6231	7.58	17894	10.34	7932	14.21

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.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG NO.: BN110724

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Nov 7 2024 12:00AM

Lab File ID: BN034887.D Time Analyzed: 14:24

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	5878	7.575	16847	10.34	7757	14.21
UPPER LIMIT	11756	8.075	33694	10.84	15514	14.711
LOWER LIMIT	2939	7.075	8423.5	9.84	3878.5	13.711
EPA SAMPLE NO.						
01 ICVBN110724	5596	7.58	16947	10.34	8617	14.21
02 PB164229BL	6231	7.58	17894	10.34	7932	14.21

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: CHEMTECH

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG NO.: BN110724

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Nov 7 2024 12:00AM

Lab File ID: BN034887.D Time Analyzed: 09:18

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	14504	16.957	8221	21.151	6835	23.317
UPPER LIMIT	29008	17.457	16442	21.651	13670	23.817
LOWER LIMIT	7252	16.457	4110.5	20.651	3417.5	22.817
EPA SAMPLE NO.						
01 SSTDCCC0.4	15173	16.95	8034	21.15	6346	23.32
02 ICVBN110724	17630	16.95	11730	21.15	10292	23.32
03 PB164229BL	16492	16.95	9359	21.15	7338	23.32

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.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN110724 SAS No.: BN110724 SDG NO.: BN110724

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Nov 7 2024 12:00AM

Lab File ID: BN034887.D Time Analyzed: 14:24

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	14504	16.957	8221	21.151	6835	23.317
UPPER LIMIT	29008	17.457	16442	21.651	13670	23.817
LOWER LIMIT	7252	16.457	4110.5	20.651	3417.5	22.817
EPA SAMPLE NO.						
01 ICVBN110724	17630	16.95	11730	21.15	10292	23.32
02 PB164229BL	16492	16.95	9359	21.15	7338	23.32

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.

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: BN110724

Client: _____

Analytical Method: **8270-Modified**

DataFile: _____

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	RPD		Limits		RPD
							Qual	Qual	Low	High	
P4368-08	n-Nitrosodimethylamine	0.2	0.22	ug/L	110				70	130	
	bis(2-Chloroethyl)ether	0.1	0.22	ug/L	220		*		70	130	
	Naphthalene	0.1	0.22	ug/L	220		*		70	130	
	Hexachlorobutadiene	0.1	0.21	ug/L	210		*		70	130	
	2-Methylnaphthalene	0.1	0.22	ug/L	220		*		70	130	
	Acenaphthylene	0.1	0.21	ug/L	210		*		70	130	
	Acenaphthene	0.1	0.2	ug/L	200		*		70	130	
	Fluorene	0.1	0.21	ug/L	210		*		70	130	
	4,6-Dinitro-2-methylphenol	0.2	0.19	ug/L	95				70	130	
	4-Bromophenyl-phenylether	0.1	0.2	ug/L	200		*		70	130	
	Hexachlorobenzene	0.1	0.22	ug/L	220		*		70	130	
	Atrazine	0.1	0.2	ug/L	200		*		70	130	
	Pentachlorophenol	0.2	0.18	ug/L	90				70	130	
	Phenanthrene	0.1	0.22	ug/L	220		*		70	130	
	Anthracene	0.1	0.22	ug/L	220		*		70	130	
	Fluoranthene	0.1	0.21	ug/L	210		*		70	130	
	Pyrene	0.1	0.21	ug/L	210		*		70	130	
	Benzo(a)anthracene	0.1	0.22	ug/L	220		*		70	130	
	Chrysene	0.1	0.22	ug/L	220		*		70	130	
	bis(2-Ethylhexyl)phthalate	0.2	0.19	ug/L	95				70	130	
	Benzo(b)fluoranthene	0.1	0.23	ug/L	230		*		70	130	
	Benzo(k)fluoranthene	0.1	0.22	ug/L	220		*		70	130	
	Benzo(a)pyrene	0.1	0.22	ug/L	220		*		70	130	
	Indeno(1,2,3-cd)pyrene	0.1	0.2	ug/L	200		*		70	130	
	Dibenz(a,h)anthracene	0.1	0.2	ug/L	200		*		70	130	
	Benzo(g,h,i)perylene	0.1	0.21	ug/L	210		*		70	130	
	1,4-Dioxane	0.2	0.23	ug/L	115				70	130	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: BN110724

Client: _____

Analytical Method: **8270-Modified**

DataFile: _____

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	RPD		Low	Limits	
							Qual	Qual		High	RPD
P4368-02	n-Nitrosodimethylamine	6.6	6.8	ug/Kg	103				70	130	
	bis(2-Chloroethyl)ether	3.3	7.3	ug/Kg	221		*		70	130	
	Naphthalene	3.3	7.2	ug/Kg	218		*		70	130	
	Hexachlorobutadiene	3.3	7	ug/Kg	212		*		70	130	
	2-Methylnaphthalene	3.3	7.1	ug/Kg	215		*		70	130	
	Acenaphthylene	3.3	7	ug/Kg	212		*		70	130	
	Acenaphthene	3.3	6.8	ug/Kg	206		*		70	130	
	Fluorene	3.3	6.8	ug/Kg	206		*		70	130	
	4,6-Dinitro-2-methylphenol	6.6	6.1	ug/Kg	92				70	130	
	4-Bromophenyl-phenylether	3.3	6.9	ug/Kg	209		*		70	130	
	Hexachlorobenzene	3.3	7.1	ug/Kg	215		*		70	130	
	Atrazine	3.3	6.7	ug/Kg	203		*		70	130	
	Pentachlorophenol	6.6	5.9	ug/Kg	89				70	130	
	Phenanthrene	3.3	7.3	ug/Kg	221		*		70	130	
	Anthracene	3.3	7.4	ug/Kg	224		*		70	130	
	Fluoranthene	3.3	7.1	ug/Kg	215		*		70	130	
	Pyrene	3.3	7.1	ug/Kg	215		*		70	130	
	Benzo(a)anthracene	3.3	7.3	ug/Kg	221		*		70	130	
	Chrysene	3.3	7.6	ug/Kg	230		*		70	130	
	bis(2-Ethylhexyl)phthalate	6.6	6.4	ug/Kg	97				70	130	
	Benzo(b)fluoranthene	3.3	8	ug/Kg	242		*		70	130	
	Benzo(k)fluoranthene	3.3	7.3	ug/Kg	221		*		70	130	
	Benzo(a)pyrene	3.3	7.1	ug/Kg	215		*		70	130	
	Indeno(1,2,3-cd)pyrene	3.3	6.5	ug/Kg	197		*		70	130	
	Dibenz(a,h)anthracene	3.3	6.6	ug/Kg	200		*		70	130	
	Benzo(g,h,i)perylene	3.3	7.1	ug/Kg	215		*		70	130	
	1,4-Dioxane	6.6	7	ug/Kg	106				70	130	

4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

SSTDCCC0.4

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BN110724

SAS No.: BN110724 SDG NO.: BN110724

Lab File ID: BN034884.D

Lab Sample ID: SSTDCCC0.4

Instrument ID: BNA_N

Date Extracted: _____

Matrix: (soil/water) Water

Date Analyzed: 11/07/2024

Level: (low/med) LOW

Time Analyzed: 09:18

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
SSTDCCC0.4	SSTDCCC0.4	BN034884.D	11/07/2024
ICVBN110724	SSTDICV0.4	BN034892.D	11/07/2024

COMMENTS: _____

4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

PB164229BL

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BN110724

SAS No.: BN110724 SDG NO.: BN110724

Lab File ID: BN034893.D

Lab Sample ID: PB164229BL

Instrument ID: BNA_N

Date Extracted: 10/17/2024

Matrix: (soil/water) Water

Date Analyzed: 11/07/2024

Level: (low/med) LOW

Time Analyzed: 15:04

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED

COMMENTS:

Surrogate Summary

SW-846

SDG No.: BN110724

Client: _____

Analytical Method: **8270-Modified**

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
SSTDCCC0.4	SSTDCCC0.4	BN034884.D	2-Methylnaphthalene-d10	0.4	362.00	90500	*	20	139
			Fluoranthene-d10	0.4	344.00	86000	*	30	150
			2-Fluorophenol	0.4	355.00	88750	*	10	100
			Phenol-d6	0.4	348.00	87000	*	10	100
			Nitrobenzene-d5	0.4	369.00	92250	*	27	123
			2-Fluorobiphenyl	0.4	367.00	91750	*	34	132
			2,4,6-Tribromophenol	0.4	326.00	81500	*	10	131
			Terphenyl-d14	0.4	405.00	101250	*	35	157
SSTDICV0.4	ICVBN110724	BN034892.D	2-Methylnaphthalene-d10	0.4	379.00	94750	*	20	139
			Fluoranthene-d10	0.4	383.00	95750	*	30	150
			2-Fluorophenol	0.4	354.00	88500	*	10	100
			Phenol-d6	0.4	348.00	87000	*	10	100
			Nitrobenzene-d5	0.4	345.00	86250	*	27	123
			2-Fluorobiphenyl	0.4	359.00	89750	*	34	132
			2,4,6-Tribromophenol	0.4	383.00	95750	*	10	131
			Terphenyl-d14	0.4	385.00	96250	*	35	157

Surrogate Summary

SW-846

SDG No.: BN110724

Client: _____

Analytical Method: **8270-Modified**

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
PB164229BL	PB164229BL	BN034893.D	2-Methylnaphthalene-d10	0.4	0.38	95		20	139
			Fluoranthene-d10	0.4	0.40	100		30	150
			2-Fluorophenol	0.4	0.36	91		10	100
			Phenol-d6	0.4	0.36	90		10	100
			Nitrobenzene-d5	0.4	0.38	95		27	123
			2-Fluorobiphenyl	0.4	0.41	101		34	132
			2,4,6-Tribromophenol	0.4	0.28	70		10	131
			Terphenyl-d14	0.4	0.52	130		35	157

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: CHEMTECH

Contract: BN110724

Lab Code: CHEM

SAS No.: BN110724 SDG NO.: BN110724

Lab File ID: BN034883.D

DFTPP Injection Date: 11/07/2024

Instrument ID: BNA_N

DFTPP Injection Time: 08:39

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	64.6
68	Less than 2.0% of mass 69	0.9 (1.6) 1
69	Mass 69 relative abundance	57.5
70	Less than 2.0% of mass 69	0.3 (0.6) 1
127	10.0 - 80.0% of mass 198	64.5
197	Less than 2.0% of mass 198	0.3
198	Base Peak, 100% relative abundance	100
199	5.0 to 9.0% of mass 198	7
275	10.0 - 60.0% of mass 198	19.4
365	Greater than 1% of mass 198	2.3
441	Present, but less than mass 443	6.6
442	Greater than 50% of mass 198	37.1
443	15.0 - 24.0% of mass 442	8.9 (24) 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	BN034884.D	11/07/2024	09:18
SSTDICC0.1	SSTDICC0.1	BN034885.D	11/07/2024	10:02
SSTDICC0.2	SSTDICC0.2	BN034886.D	11/07/2024	10:48
SSTDICCC0.4	SSTDICCC0.4	BN034887.D	11/07/2024	11:24
SSTDICC0.8	SSTDICC0.8	BN034888.D	11/07/2024	12:00
SSTDICC1.6	SSTDICC1.6	BN034889.D	11/07/2024	12:36
SSTDICC3.2	SSTDICC3.2	BN034890.D	11/07/2024	13:13
SSTDICC5.0	SSTDICC5.0	BN034891.D	11/07/2024	13:49
ICVBN110724	SSTDICV0.4	BN034892.D	11/07/2024	14:24
PB164229BL	PB164229BL	BN034893.D	11/07/2024	15:04
SSTDCCC0.4EC	SSTDCCC0.4	BN034896.D	11/07/2024	16:58