

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: _____

Lab Code: CHEM Case No.: BN012224 SAS No.: BN012224 SDG No.: BN012224

Instrument ID: BNA_N Calibration Date/Time: 1/22/2024 1:16:00

Lab File ID: BN029333.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	0.662	0.611		-7.704	20.0
Fluoranthene-d10	1.185	1.070		-9.705	20.0
n-Nitrosodimethylamine	0.527	0.477		-9.488	20.0
2-Fluorophenol	1.062	0.982		-7.533	20.0
Phenol-d6	1.419	1.319		-7.047	20.0
bis(2-Chloroethyl)ether	1.310	1.240		-5.344	20.0
Nitrobenzene-d5	0.424	0.386		-8.962	20.0
Naphthalene	1.286	1.182		-8.087	20.0
Hexachlorobutadiene	0.271	0.258		-4.797	20.0
2-Methylnaphthalene	0.836	0.766		-8.373	20.0
2-Fluorobiphenyl	1.926	1.781		-7.529	20.0
Acenaphthylene	2.133	1.910		-10.455	20.0
Acenaphthene	1.434	1.295		-9.693	20.0
Fluorene	1.945	1.776		-8.689	20.0
4,6-Dinitro-2-methylphenol	0.074	0.065		-12.162	20.0
2,4,6-Tribromophenol	0.208	0.180		-13.462	20.0
4-Bromophenyl-phenylether	0.292	0.268		-8.219	20.0
Hexachlorobenzene	0.360	0.336		-6.667	20.0
Atrazine	0.214	0.188		-12.15	20.0
Pentachlorophenol	0.121	0.103		-14.876	20.0
Phenanthrene	1.372	1.254		-8.601	20.0
Anthracene	1.213	1.065		-12.201	20.0
Fluoranthene	1.643	1.482		-9.799	20.0
Pyrene	1.824	1.700		-6.798	20.0
Terphenyl-d14	1.025	0.946		-7.707	20.0
Benzo(a)anthracene	1.668	1.482		-11.151	20.0
Chrysene	1.719	1.595		-7.213	20.0
Bis(2-ethylhexyl)phthalate	0.919	0.763		-16.975	20.0
Benzo(b)fluoranthene	1.703	1.572		-7.692	20.0
Benzo(k)fluoranthene	1.672	1.537		-8.074	20.0
Benzo(a)pyrene	1.340	1.179		-12.015	20.0
Indeno(1,2,3-cd)pyrene	1.709	1.504		-11.995	20.0
Dibenzo(a,h)anthracene	1.382	1.225		-11.36	20.0
Benzo(g,h,i)perylene	1.482	1.328		-10.391	20.0
1,4-Dioxane	0.520	0.494		-5.0	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:	SSTDICV0.4	SDG No.:	BN012224
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
BN029338.D	1		1/22/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	360		29	200	200	ug/L
111-44-4	bis(2-Chloroethyl)ether	380		21	100	100	ug/L
91-20-3	Naphthalene	370		19	100	100	ug/L
87-68-3	Hexachlorobutadiene	380		19	100	100	ug/L
91-57-6	2-Methylnaphthalene	370		21	100	100	ug/L
208-96-8	Acenaphthylene	360		18	100	100	ug/L
83-32-9	Acenaphthene	370		16	100	100	ug/L
Total PAH	Total PAH	5660		400	100	1600	ug/L
86-73-7	Fluorene	360		17	100	100	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	320		73	200	200	ug/L
101-55-3	4-Bromophenyl-phenylether	340		20	100	100	ug/L
118-74-1	Hexachlorobenzene	370		24	100	100	ug/L
1912-24-9	Atrazine	340		18	100	100	ug/L
87-86-5	Pentachlorophenol	310		59	200	200	ug/L
85-01-8	Phenanthrene	360		19	100	100	ug/L
120-12-7	Anthracene	350		22	100	100	ug/L
206-44-0	Fluoranthene	350		19	100	100	ug/L
129-00-0	Pyrene	360		26	100	100	ug/L
56-55-3	Benzo(a)anthracene	360		24	100	100	ug/L
218-01-9	Chrysene	370		20	100	100	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	320		81	200	200	ug/L
205-99-2	Benzo(b)fluoranthene	340		24	100	100	ug/L
207-08-9	Benzo(k)fluoranthene	350		27	100	100	ug/L
50-32-8	Benzo(a)pyrene	340		36	100	100	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	340		40	100	100	ug/L
53-70-3	Dibenzo(a,h)anthracene	340		42	100	100	ug/L
191-24-2	Benzo(g,h,i)perylene	340		31	100	100	ug/L
123-91-1	1,4-Dioxane	400		61	200	200	ug/L
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	369	*	30-150		92250	SPK: -0.4
93951-69-0	Fluoranthene-d10	353	*	30-150		88250	SPK: -0.4

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	SSTDICV0.4			SDG No.:	BN012224		
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
BN029338.D	1		1/22/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
367-12-4	2-Fluorophenol	370	*	10-92		92500	SPK: -0.4
13127-88-3	Phenol-d6	368	*	10-100		92000	SPK: -0.4
4165-60-0	Nitrobenzene-d5	353	*	11-175		88250	SPK: -0.4
321-60-8	2-Fluorobiphenyl	374	*	10-175		93500	SPK: -0.4
118-79-6	2,4,6-Tribromophenol	299	*	24-148		74750	SPK: -0.4
1718-51-0	Terphenyl-d14	363	*	54-171		90750	SPK: -0.4
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	4148	7.919				
1146-65-2	Naphthalene-d8	11226	10.712				
15067-26-2	Acenaphthene-d10	6051	14.554				
1517-22-2	Phenanthrene-d10	13991	17.305				
1719-03-5	Chrysene-d12	12478	21.51				
1520-96-3	Perylene-d12	13310	23.905				

Report of Analysis

Client:		Date Collected:	1/22/2024 12:00:00 AM
Project:		Date Received:	1/22/2024 12:00:00 AM
Client Sample ID:	PB158557BL	SDG No.:	BN012224
Lab Sample ID:	PB158557BL	Matrix:	Solid
Analytical Method:	8270-Modified	% Moisture:	0
Sample Wt/Vol:	30.03 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
BN029339.D	1	1/22/2024 12:00:00 AM	1/22/2024 12:00:00 AM	PB158557

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	1.7	U	1.7	6.6	6.6	ug/Kg
111-44-4	bis(2-Chloroethyl)ether	0.79	U	0.79	3.3	3.3	ug/Kg
91-20-3	Naphthalene	0.72	U	0.72	3.3	3.3	ug/Kg
87-68-3	Hexachlorobutadiene	0.43	U	0.43	3.3	3.3	ug/Kg
91-57-6	2-Methylnaphthalene	0.8	U	0.8	3.3	3.3	ug/Kg
208-96-8	Acenaphthylene	0.61	U	0.61	3.3	3.3	ug/Kg
83-32-9	Acenaphthene	0.53	U	0.53	3.3	3.3	ug/Kg
Total PAH	Total PAH	12.81	U	12.81	3.3	52.8	ug/Kg
86-73-7	Fluorene	0.53	U	0.53	3.3	3.3	ug/Kg
534-52-1	4,6-Dinitro-2-methylphenol	1.1	U	1.1	6.6	6.6	ug/Kg
101-55-3	4-Bromophenyl-phenylether	0.74	U	0.74	3.3	3.3	ug/Kg
118-74-1	Hexachlorobenzene	0.85	U	0.85	3.3	3.3	ug/Kg
1912-24-9	Atrazine	0.49	U	0.49	3.3	3.3	ug/Kg
87-86-5	Pentachlorophenol	1.6	U	1.6	6.6	6.6	ug/Kg
85-01-8	Phenanthrene	0.63	U	0.63	3.3	3.3	ug/Kg
120-12-7	Anthracene	0.69	U	0.69	3.3	3.3	ug/Kg
206-44-0	Fluoranthene	0.56	U	0.56	3.3	3.3	ug/Kg
129-00-0	Pyrene	1.2	U	1.2	3.3	3.3	ug/Kg
56-55-3	Benzo(a)anthracene	0.76	U	0.76	3.3	3.3	ug/Kg
218-01-9	Chrysene	0.72	U	0.72	3.3	3.3	ug/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	1.5	U	1.5	6.6	6.6	ug/Kg
205-99-2	Benzo(b)fluoranthene	0.88	U	0.88	3.3	3.3	ug/Kg
207-08-9	Benzo(k)fluoranthene	0.73	U	0.73	3.3	3.3	ug/Kg
50-32-8	Benzo(a)pyrene	0.75	U	0.75	3.3	3.3	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	1.4	U	1.4	3.3	3.3	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	1.1	U	1.1	3.3	3.3	ug/Kg
191-24-2	Benzo(g,h,i)perylene	1	U	1	3.3	3.3	ug/Kg
123-91-1	1,4-Dioxane	1.7	U	1.7	6.6	6.6	ug/Kg
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	0.276		21-128		69	SPK: -0.4
93951-69-0	Fluoranthene-d10	0.273		13-127		68	SPK: -0.4

**Report of Analysis**

Client:		Date Collected:	1/22/2024 12:00:00 AM
Project:		Date Received:	1/22/2024 12:00:00 AM
Client Sample ID:	PB158557BL	SDG No.:	BN012224
Lab Sample ID:	PB158557BL	Matrix:	Solid
Analytical Method:	8270-Modified	% Moisture:	0
Sample Wt/Vol:	30.03	Units:	g
Soil Aliquot Vol:		Final Vol:	1000 uL
Extraction Type :		Test:	
	Decanted :	Level :	LOW
Injection Volume :		GPC Factor :	
		GPC Cleanup :	PH :

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BN029339.D	1	1/22/2024 12:00:00 AM	1/22/2024 12:00:00 AM	PB158557

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
367-12-4	2-Fluorophenol	0.351		28-122		88	SPK: -0.4
13127-88-3	Phenol-d6	0.335		12-126		84	SPK: -0.4
4165-60-0	Nitrobenzene-d5	0.276		14-112		69	SPK: -0.4
321-60-8	2-Fluorobiphenyl	0.314		22-114		79	SPK: -0.4
118-79-6	2,4,6-Tribromophenol	0.207		10-137		52	SPK: -0.4
1718-51-0	Terphenyl-d14	0.302		21-130		75	SPK: -0.4
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	4528	7.919				
1146-65-2	Naphthalene-d8	11841	10.712				
15067-26-2	Acenaphthene-d10	5736	14.554				
1517-22-2	Phenanthrene-d10	13831	17.305				
1719-03-5	Chrysene-d12	11514	21.51				
1520-96-3	Perylene-d12	11511	23.908				



Hit Summary Sheet
SW-846

SDG No.: BN012224

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
Client ID :	SSTDICV0.4							
SSTDICV0.4	SSTDICV0.4	Water	1,4-Dioxane	400.000	61		200	ug/L
SSTDICV0.4	SSTDICV0.4	Water	2-Methylnaphthalene	370.000	21		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	4,6-Dinitro-2-methylphenol	320.000	73		200	ug/L
SSTDICV0.4	SSTDICV0.4	Water	4-Bromophenyl-phenylether	340.000	20		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Acenaphthene	370.000	16		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Acenaphthylene	360.000	18		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Anthracene	350.000	22		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Atrazine	340.000	18		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Benzo(a)anthracene	360.000	24		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Benzo(a)pyrene	340.000	36		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Benzo(b)fluoranthene	340.000	24		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Benzo(g,h,i)perylene	340.000	31		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Benzo(k)fluoranthene	350.000	27		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	bis(2-Chloroethyl)ether	380.000	21		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Bis(2-ethylhexyl)phthalate	320.000	81		200	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Chrysene	370.000	20		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Dibenzo(a,h)anthracene	340.000	42		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Fluoranthene	350.000	19		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Fluorene	360.000	17		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Hexachlorobenzene	370.000	24		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Hexachlorobutadiene	380.000	19		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Indeno(1,2,3-cd)pyrene	340.000	40		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	n-Nitrosodimethylamine	360.000	29		200	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Naphthalene	370.000	19		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Pentachlorophenol	310.000	59		200	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Phenanthrene	360.000	19		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Pyrene	360.000	26		100	ug/L
SSTDICV0.4	SSTDICV0.4	Water	Total PAH	5,660.000	400		1600	ug/L
Total Svoc :				15,210.00				
Total Concentration:				15,210.00				



6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH

Contract:

Lab Code: CHEM Case No.: BN012224

SAS No.: BN012224

SDG No.: BN012224

Instrument ID:

Calibration Date(s): 1/22/2024 :

Calibration Time(s): 14:48

LAB FILE ID:		RRF0.1 = BN029331.D			RRF0.2 = BN029332.D		RRF0.4 = BN029333.D		
		RRF0.8 = BN029334.D			RRF1.6 = BN029335.D		RRF3.2 = BN029336.D		
COMPOUND	RRF0.1	RRF0.2	RRF0.4	RRF0.8	RRF1.6	RRF3.2	RRF	% RSD	
2-Methylnaphthalene-d10	0.633	0.631	0.611	0.684	0.715	0.703	0.662	5.9	
Fluoranthene-d10	1.110	1.102	1.070	1.192	1.286	1.310	1.185	8.0	
n-Nitrosodimethylamine	0.452	0.464	0.477	0.555	0.597	0.596	0.527	11.8	
2-Fluorophenol	1.139	1.052	0.982	1.076	1.101	1.075	1.062	5.1	
Phenol-d6	1.549	1.394	1.319	1.407	1.466	1.437	1.419	5.2	
bis(2-Chloroethyl)ether	1.353	1.290	1.240	1.342	1.376	1.340	1.310	4.4	
Nitrobenzene-d5	0.479	0.432	0.386	0.425	0.431	0.419	0.424	7.1	
Naphthalene	1.341	1.284	1.182	1.327	1.344	1.310	1.286	5.1	
Hexachlorobutadiene	0.272	0.276	0.258	0.287	0.288	0.270	0.271	5.8	
2-Methylnaphthalene	0.844	0.812	0.766	0.860	0.889	0.872	0.836	5.1	
2-Fluorobiphenyl	2.071	1.968	1.781	1.977	2.019	1.905	1.926	6.1	
Acenaphthylene	2.043	1.985	1.910	2.163	2.316	2.305	2.133	7.4	
Acenaphthene	1.501	1.405	1.295	1.454	1.523	1.480	1.434	5.6	
Fluorene	2.017	1.916	1.776	1.971	2.069	2.004	1.945	5.2	
4,6-Dinitro-2-methylphenol	0.067	0.062	0.065	0.073	0.080	0.087	0.074	12.9	
2,4,6-Tribromophenol	0.243	0.206	0.180	0.194	0.210	0.218	0.208	9.5	
4-Bromophenyl-phenylether	0.305	0.294	0.268	0.293	0.301	0.302	0.292	4.5	
Hexachlorobenzene	0.367	0.365	0.336	0.370	0.375	0.366	0.360	4.0	
Atrazine	0.196	0.198	0.188	0.212	0.234	0.239	0.214	9.5	
Pentachlorophenol	0.094	0.098	0.103	0.120	0.135	0.148	0.121	18.9	
Phenanthrene	1.431	1.368	1.254	1.368	1.418	1.423	1.372	4.5	
Anthracene	1.191	1.140	1.065	1.194	1.291	1.338	1.213	7.8	
Fluoranthene	1.582	1.530	1.482	1.654	1.786	1.805	1.643	7.4	
Pyrene	1.973	1.846	1.700	1.817	1.890	1.815	1.824	5.1	
Terphenyl-d14	1.147	1.045	0.946	1.028	1.054	1.009	1.025	6.7	
Benzo (a) anthracene	1.646	1.580	1.482	1.667	1.824	1.797	1.668	7.1	
Chrysene	1.753	1.691	1.595	1.757	1.852	1.758	1.719	5.1	
Bis (2-ethylhexyl)phthalate	1.280	0.866	0.763	0.803	0.864	0.949	0.919	18.6	
Benzo (b) fluoranthene	1.732	1.644	1.572	1.752	1.794	1.736	1.703	4.4	
Benzo (k) fluoranthene	1.612	1.568	1.537	1.765	1.787	1.759	1.672	6.1	
Benzo (a) pyrene	1.249	1.218	1.179	1.363	1.460	1.468	1.340	9.2	
Indeno (1,2,3-cd) pyrene	1.504	1.504	1.504	1.724	1.937	1.915	1.709	11.9	
Dibenzo (a,h) anthracene	1.191	1.198	1.225	1.413	1.583	1.551	1.382	12.6	
Benzo (g,h,i) perylene	1.377	1.348	1.328	1.486	1.652	1.607	1.482	9.0	
1,4-Dioxane	0.553	0.499	0.494	0.531	0.552	0.523	0.520	5.3	

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECHLab Code: CHEM Case No.: BN012224 SAS No.: BN012224 SDG NO.: BN012224EPA Sample No.: SSTDICCC0.4 Date Analyzed: Jan 22 2024 12:00AMLab File ID: BN029333.D Time Analyzed: 19:36Instrument 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	3548	7.919	9670	10.71	5316	14.55
UPPER LIMIT	7096	8.419	19340	11.211	10632	15.053
LOWER LIMIT	1774	7.419	4835	10.211	2658	14.053
EPA SAMPLE NO.						
01 SSTDICV0.4	4148	7.92	11226	10.71	6051	14.55
02 PB158557BL	4528	7.92	11841	10.71	5736	14.55

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.

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN012224 SAS No.: BN012224 SDG NO.: BN012224

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Jan 22 2024 12:00AM

Lab File ID: BN029333.D Time Analyzed: 19:36

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	3548	7.919	9670	10.71	5316	14.55
UPPER LIMIT	7096	8.419	19340	11.211	10632	15.053
LOWER LIMIT	1774	7.419	4835	10.211	2658	14.053
EPA SAMPLE NO.						
01 SSTDICV0.4	4148	7.92	11226	10.71	6051	14.55
02 PB158557BL	4528	7.92	11841	10.71	5736	14.55

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.

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN012224 SAS No.: BN012224 SDG NO.: BN012224

EPA Sample No.: SSTDICC0.4 Date Analyzed: Jan 22 2024 12:00AM

Lab File ID: BN029333.D Time Analyzed: 19:36

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	12265	17.305	10934	21.51	10903	23.905
UPPER LIMIT	24530	17.805	21868	22.01	21806	24.405
LOWER LIMIT	6132.5	16.805	5467	21.01	5451.5	23.405
EPA SAMPLE NO.						
01 SSTDICV0.4	13991	17.31	12478	21.51	13310	23.91
02 PB158557BL	13831	17.31	11514	21.51	11511	23.91

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.

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BN012224 SAS No.: BN012224 SDG NO.: BN012224

EPA Sample No.: SSTDICCC0.4 Date Analyzed: Jan 22 2024 12:00AM

Lab File ID: BN029333.D Time Analyzed: 19:36

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	12265	17.305	10934	21.51	10903	23.905
UPPER LIMIT	24530	17.805	21868	22.01	21806	24.405
LOWER LIMIT	6132.5	16.805	5467	21.01	5451.5	23.405
EPA SAMPLE NO.						
01 SSTDICV0.4	13991	17.31	12478	21.51	13310	23.91
02 PB158557BL	13831	17.31	11514	21.51	11511	23.91

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.



4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

SSTDICV0.4

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BN012224SAS No.: BN012224 SDG NO.: BN012224Lab File ID: BN029338.DLab Sample ID: SSTDICV0.4Instrument ID: BNA_N

Date Extracted: _____

Matrix: (soil/water) WaterDate Analyzed: 01/22/2024Level: (low/med) LOWTime Analyzed: 19:36

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
SSTDICV0.4	SSTDICV0.4	BN029338.D	01/22/2024

COMMENTS: _____



4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

PB158557BL

Lab Name: CHEMTECH

Contract:

Lab Code: CHEM Case No.: BN012224

SAS No.: BN012224 SDG NO.: BN012224

Lab File ID: BN029339.D

Lab Sample ID: PB158557BL

Instrument ID: BNA_N

Date Extracted: 01/22/2024

Matrix: (soil/water) Solid

Date Analyzed: 01/22/2024

Level: (low/med) LOW

Time Analyzed: 20:12

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.: BN012224

Client:

Analytical Method: 8270-Modified

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
SSTDICV0.4	SSTDICV0.4	BN029338.D	2-Methylnaphthalene-d10	0.4	369.00	92250	*	30	150
			Fluoranthene-d10	0.4	353.00	88250	*	30	150
			2-Fluorophenol	0.4	370.00	92500	*	10	92
			Phenol-d6	0.4	368.00	92000	*	10	100
			Nitrobenzene-d5	0.4	353.00	88250	*	11	175
			2-Fluorobiphenyl	0.4	374.00	93500	*	10	175
			2,4,6-Tribromophenol	0.4	299.00	74750	*	24	148
			Terphenyl-d14	0.4	363.00	90750	*	54	171



Surrogate Summary
SW-846

SDG No.: BN012224

Client:

Analytical Method: 8270-Modified

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
PB158557BL	PB158557BL	BN029339.D	2-Methylnaphthalene-d10	0.4	0.28	69		21	128
			Fluoranthene-d10	0.4	0.27	68		13	127
			2-Fluorophenol	0.4	0.35	88		28	122
			Phenol-d6	0.4	0.34	84		12	126
			Nitrobenzene-d5	0.4	0.28	69		14	112
			2-Fluorobiphenyl	0.4	0.31	79		22	114
			2,4,6-Tribromophenol	0.4	0.21	52		10	137
			Terphenyl-d14	0.4	0.30	75		21	130



5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)Lab Name: CHEMTECHContract: BN012224Lab Code: CHEMSAS No.: BN012224 SDG NO.: BN012224Lab File ID: BN029330.DDFTPP Injection Date: 01/22/2024Instrument ID: BNA_NDFTPP Injection Time: 14:12

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	45.1
68	Less than 2.0% of mass 69	0.9 (1.7) 1
69	Mass 69 relative abundance	52.5
70	Less than 2.0% of mass 69	0.3 (0.6) 1
127	10.0 - 80.0% of mass 198	55.8
197	Less than 2.0% of mass 198	0.8
198	Base Peak, 100% relative abundance	100
199	5.0 to 9.0% of mass 198	6.7
275	10.0 - 60.0% of mass 198	26.2
365	Greater than 1% of mass 198	3.4
441	Present, but less than mass 443	12.1
442	Greater than 50% of mass 198	78.1
443	15.0 - 24.0% of mass 442	12.6 (16.1) 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	BN029331.D	01/22/2024	14:48
SSTDICC0.2	SSTDICC0.2	BN029332.D	01/22/2024	15:24
SSTDICCC0.4	SSTDICCC0.4	BN029333.D	01/22/2024	16:00
SSTDICC0.8	SSTDICC0.8	BN029334.D	01/22/2024	16:36
SSTDICC1.6	SSTDICC1.6	BN029335.D	01/22/2024	17:12
SSTDICC3.2	SSTDICC3.2	BN029336.D	01/22/2024	17:48
SSTDICC5.0	SSTDICC5.0	BN029337.D	01/22/2024	18:24
SSTDICV0.4	SSTDICV0.4	BN029338.D	01/22/2024	19:36
PB158557BL	PB158557BL	BN029339.D	01/22/2024	20:12