

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: _____

Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG No.: BP091624

Instrument ID: BNA_P Calibration Date/Time: 9/16/2024 1:13:53

Lab File ID: BP021932.D Init. Calib. Date(s): _____

EPA Sample No.: SSTDICCC040 Init. Calib. Time(s): _____

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

COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
n-Nitrosodimethylamine	0.669	0.681		1.794	
Pyridine	1.117	1.248		11.728	
2-Fluorophenol	1.108	1.058		-4.513	
Benzaldehyde	0.724	0.771		6.492	
Aniline	1.227	1.246		1.548	
Phenol-d6	1.358	1.390		2.356	
Phenol	1.352	1.385		2.441	20.0
bis(2-Chloroethyl) ether	1.084	1.106		2.03	
2-Chlorophenol	1.155	1.161		0.519	
1,2-Dichlorobenzene	1.392	1.391		-0.072	
1,3-Dichlorobenzene	1.437	1.450		0.905	
1,4-Dichlorobenzene	1.455	1.455		0.0	20.0
Benzyl Alcohol	1.042	1.026		-1.536	
2-Methylphenol	0.944	0.962		1.907	
2,2-oxybis(1-Chloropropane)	1.200	1.268		5.667	
Acetophenone	0.470	0.359		-23.617	
3+4-Methylphenols	1.324	1.357		2.492	
n-Nitroso-di-n-propylamine	0.913	0.944	0.050	3.395	
Nitrobenzene-d5	0.365	0.281		-23.014	
Hexachloroethane	0.535	0.557		4.112	
Nitrobenzene	0.377	0.292		-22.546	
Isophorone	0.583	0.486		-16.638	
2-Nitrophenol	0.149	0.122		-18.121	20.0
2,4-Dimethylphenol	0.185	0.153		-17.297	
bis(2-Chloroethoxy) methane	0.386	0.397		2.85	
2,4-Dichlorophenol	0.316	0.321		1.582	20.0
1,2,4-Trichlorobenzene	0.385	0.384		-0.26	
Benzoic acid	0.202	0.165		-18.317	
Naphthalene	1.000	1.002		0.2	
4-Chloroaniline	0.368	0.372		1.087	
Hexachlorobutadiene	0.266	0.253		-4.887	20.0
Caprolactam	0.096	0.101		5.208	
4-Chloro-3-methylphenol	0.343	0.338		-1.458	20.0
2-Methylnaphthalene	0.727	0.718		-1.238	
Hexachlorocyclopentadiene	0.209	0.199	0.050	-4.785	
2,4,6-Trichlorophenol	0.384	0.390		1.563	20.0
2-Fluorobiphenyl	1.312	1.268		-3.354	
2,4,5-Trichlorophenol	0.440	0.439		-0.227	
1,1-Biphenyl	1.363	1.339		-1.761	

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Instrument ID: BNA_P Calibration Date/Time: 9/16/2024 1:13:53

Lab File ID: BP021932.D Init. Calib. Date(s): _____

EPA Sample No.: SSTDICCC040 Init. Calib. Time(s): _____

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

COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
2-Chloronaphthalene	1.094	1.081		-1.188	
2-Nitroaniline	0.301	0.309		2.658	
Dimethylphthalate	1.432	1.392		-2.793	
Acenaphthylene	1.567	1.566		-0.064	
2,6-Dinitrotoluene	0.315	0.319		1.27	
3-Nitroaniline	0.289	0.302		4.498	
Acenaphthene	1.021	1.002		-1.861	20.0
2,4-Dinitrophenol	0.164	0.164	0.050	0.0	
4-Nitrophenol	0.252	0.268	0.050	6.349	
Dibenzofuran	1.631	1.674		2.636	
2,4-Dinitrotoluene	0.426	0.444		4.225	
Diethylphthalate	1.414	1.393		-1.485	
4-Chlorophenyl-phenylether	0.716	0.725		1.257	
Fluorene	1.335	1.365		2.247	
4-Nitroaniline	0.305	0.322		5.574	
4,6-Dinitro-2-methylphenol	0.106	0.106		0.0	
n-Nitrosodiphenylamine	0.512	0.506		-1.172	20.0
Azobenzene	1.234	1.273		3.16	
2,4,6-Tribromophenol	0.263	0.254		-3.422	
4-Bromophenyl-phenylether	0.204	0.199		-2.451	
Hexachlorobenzene	0.243	0.232		-4.527	
Atrazine	0.187	0.185		-1.07	
Pentachlorophenol	0.163	0.160		-1.84	20.0
Phenanthrene	0.983	0.974		-0.916	
Anthracene	0.944	0.941		-0.318	
Carbazole	0.936	0.939		0.321	
Di-n-butylphthalate	1.092	1.073		-1.74	
Fluoranthene	1.283	1.290		0.546	20.0
Benzidine	0.291	0.228		-21.649	
Pyrene	1.278	1.312		2.66	
Terphenyl-d14	1.013	0.983		-2.962	
Butylbenzylphthalate	0.483	0.482		-0.207	
3,3-Dichlorobenzidine	0.429	0.441		2.797	
Benzo(a)anthracene	1.288	1.276		-0.932	
Chrysene	1.232	1.221		-0.893	
Bis(2-ethylhexyl)phthalate	0.705	0.711		0.851	
Di-n-octyl phthalate	1.163	1.172		0.774	20.0
Benzo(b)fluoranthene	1.143	1.179		3.15	
Benzo(k)fluoranthene	1.085	1.138		4.885	

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Lab Name: CHEMTECH Contract: _____

Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG No.: BP091624

Instrument ID: BNA_P Calibration Date/Time: 9/16/2024 1:13:53

Lab File ID: BP021932.D Init. Calib. Date(s): _____

EPA Sample No.: SSTDICCC040 Init. Calib. Time(s): _____

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

COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
Benzo(a)pyrene	1.016	1.017		0.098	20.0
Indeno(1,2,3-cd)pyrene	1.284	1.258		-2.025	
Dibenzo(a,h)anthracene	1.063	1.031		-3.01	
Benzo(g,h,i)perylene	1.106	1.090		-1.447	
1,2,4,5-Tetrachlorobenzene	0.600	0.590		-1.667	
1,4-Dioxane	0.481	0.615		27.859	20.0
2,3,4,6-Tetrachlorophenol	0.375	0.391		4.267	
1-Methylnaphthalene	0.721	0.715		-0.832	

All other compounds must meet a minimum RRF of 0.010.

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBP091624			SDG No.:	BP091624		
Lab Sample ID:	SSTDICV040			Matrix:	Water		
Analytical Method:	8270E			% 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
BP021936.D	1		9/16/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	36000		1000	8000	10000	ug/L
110-86-1	Pyridine	40200		1600	4000	5000	ug/L
100-52-7	Benzaldehyde	43400		4000	8000	10000	ug/L
62-53-3	Aniline	38300		1600	4000	5000	ug/L
108-95-2	Phenol	38900		930	4000	5000	ug/L
111-44-4	bis(2-Chloroethyl)ether	38700		1200	4000	5000	ug/L
95-57-8	2-Chlorophenol	39000		710	4000	5000	ug/L
95-50-1	1,2-Dichlorobenzene	39200		880	4000	5000	ug/L
541-73-1	1,3-Dichlorobenzene	39800		800	4000	5000	ug/L
106-46-7	1,4-Dichlorobenzene	39400		840	4000	5000	ug/L
100-51-6	Benzyl Alcohol	38800		1600	8000	10000	ug/L
95-48-7	2-Methylphenol	38800		1100	4000	5000	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	38200		1400	4000	5000	ug/L
98-86-2	Acetophenone	40200		1100	4000	5000	ug/L
65794-96-9	3+4-Methylphenols	38300		1200	8000	10000	ug/L
621-64-7	n-Nitroso-di-n-propylamine	37900		1500	2500	2500	ug/L
67-72-1	Hexachloroethane	40400		1000	4000	5000	ug/L
98-95-3	Nitrobenzene	41000		1300	4000	5000	ug/L
78-59-1	Isophorone	42600		1100	4000	5000	ug/L
88-75-5	2-Nitrophenol	44400		2000	4000	5000	ug/L
105-67-9	2,4-Dimethylphenol	42600		1500	4000	5000	ug/L
111-91-1	bis(2-Chloroethoxy)methane	39500		1000	4000	5000	ug/L
120-83-2	2,4-Dichlorophenol	40900		880	4000	5000	ug/L
120-82-1	1,2,4-Trichlorobenzene	39600		1100	4000	5000	ug/L
65-85-0	Benzoic acid	42500		5500	8000	10000	ug/L
91-20-3	Naphthalene	39200		1000	4000	5000	ug/L
106-47-8	4-Chloroaniline	38800		1300	4000	5000	ug/L
87-68-3	Hexachlorobutadiene	40700		1300	4000	5000	ug/L

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBP091624			SDG No.:	BP091624		
Lab Sample ID:	SSTDICV040			Matrix:	Water		
Analytical Method:	8270E			% 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
BP021936.D	1		9/16/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
105-60-2	Caprolactam	39800		1700	8000	10000	ug/L
59-50-7	4-Chloro-3-methylphenol	40700		840	4000	5000	ug/L
91-57-6	2-Methylnaphthalene	40000		1100	4000	5000	ug/L
77-47-4	Hexachlorocyclopentadiene	30400		5000	8000	10000	ug/L
88-06-2	2,4,6-Trichlorophenol	31800		890	4000	5000	ug/L
95-95-4	2,4,5-Trichlorophenol	31000		1000	4000	5000	ug/L
92-52-4	1,1-Biphenyl	30700		910	4000	5000	ug/L
91-58-7	2-Chloronaphthalene	30900		970	4000	5000	ug/L
88-74-4	2-Nitroaniline	32100		1400	4000	5000	ug/L
131-11-3	Dimethylphthalate	33300		930	4000	5000	ug/L
208-96-8	Acenaphthylene	40000		1000	4000	5000	ug/L
606-20-2	2,6-Dinitrotoluene	35800		1200	4000	5000	ug/L
99-09-2	3-Nitroaniline	40700		1400	4000	5000	ug/L
83-32-9	Acenaphthene	38600		810	4000	5000	ug/L
Total PAH	Total PAH	635700		17360	4000	80000	ug/L
51-28-5	2,4-Dinitrophenol	40200		6400	8000	10000	ug/L
100-02-7	4-Nitrophenol	41100		2000	8000	10000	ug/L
132-64-9	Dibenzofuran	40900		930	4000	5000	ug/L
121-14-2	2,4-Dinitrotoluene	41200		1500	4000	5000	ug/L
25321-14-6	2,4-Dinitrotoluene/2,6-Dinitrotoluene (mixture)	77000		2700	4000	10000	ug/L
84-66-2	Diethylphthalate	37900		1000	4000	5000	ug/L
7005-72-3	4-Chlorophenyl-phenylether	39700		980	4000	5000	ug/L
86-73-7	Fluorene	40200		960	4000	5000	ug/L
100-01-6	4-Nitroaniline	40500		2000	4000	5000	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	40500		3100	8000	10000	ug/L
86-30-6	n-Nitrosodiphenylamine	40600		890	4000	5000	ug/L
103-33-3	Azobenzene	39300		1200	4000	5000	ug/L
101-55-3	4-Bromophenyl-phenylether	39800		950	4000	5000	ug/L
118-74-1	Hexachlorobenzene	39400		1100	4000	5000	ug/L
1912-24-9	Atrazine	38600		1300	4000	5000	ug/L

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBP091624			SDG No.:	BP091624		
Lab Sample ID:	SSTDICV040			Matrix:	Water		
Analytical Method:	8270E			% 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
BP021936.D	1		9/16/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
87-86-5	Pentachlorophenol	40300		1900	8000	10000	ug/L
85-01-8	Phenanthrene	39700		890	4000	5000	ug/L
120-12-7	Anthracene	39800		1100	4000	5000	ug/L
86-74-8	Carbazole	40300		1200	4000	5000	ug/L
84-74-2	Di-n-butylphthalate	34900		1500	4000	5000	ug/L
206-44-0	Fluoranthene	39500		1300	4000	5000	ug/L
92-87-5	Benzidine	34400		4100	8000	10000	ug/L
129-00-0	Pyrene	40900		1100	4000	5000	ug/L
85-68-7	Butylbenzylphthalate	39200		2100	4000	5000	ug/L
91-94-1	3,3-Dichlorobenzidine	41700		1300	8000	10000	ug/L
56-55-3	Benzo(a)anthracene	39500		940	4000	5000	ug/L
218-01-9	Chrysene	38900		860	4000	5000	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	39600		1900	4000	5000	ug/L
117-84-0	Di-n-octyl phthalate	32200		2500	8000	10000	ug/L
205-99-2	Benzo(b)fluoranthene	40400		1100	4000	5000	ug/L
207-08-9	Benzo(k)fluoranthene	41700		1200	4000	5000	ug/L
50-32-8	Benzo(a)pyrene	39700		1700	4000	5000	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	39300		1000	4000	5000	ug/L
53-70-3	Dibenzo(a,h)anthracene	38900		1200	4000	5000	ug/L
191-24-2	Benzo(g,h,i)perylene	39400		1200	4000	5000	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	29100		1100	4000	5000	ug/L
123-91-1	1,4-Dioxane	36400		1300	4000	5000	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	41300		790	4000	5000	ug/L
90-12-0	1-Methylnaphthalene	40100		860	4000	5000	ug/L
SURROGATES							
367-12-4	2-Fluorophenol	75094	*	10-139		50063	SPK: -150
13127-88-3	Phenol-d6	78812	*	10-134		52541	SPK: -150
4165-60-0	Nitrobenzene-d5	82446	*	49-133		82446	SPK: -100
321-60-8	2-Fluorobiphenyl	60600	*	52-132		60600	SPK: -100

Report of Analysis

Client:				Date Collected:			
Project:				Date Received:			
Client Sample ID:	ICVBP091624			SDG No.:	BP091624		
Lab Sample ID:	SSTDICV040			Matrix:	Water		
Analytical Method:	8270E			% 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
BP021936.D	1		9/16/2024 12:00:00 AM	

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
118-79-6	2,4,6-Tribromophenol	76181	*	44-137		50787	SPK: -150
1718-51-0	Terphenyl-d14	78206	*	48-125		78206	SPK: -100
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	120317	7.746				
1146-65-2	Naphthalene-d8	444569	10.51				
15067-26-2	Acenaphthene-d10	465447	14.363				
1517-22-2	Phenanthrene-d10	1025609	17.145				
1719-03-5	Chrysene-d12	1049686	21.586				
1520-96-3	Perylene-d12	1189524	24.892				

Report of Analysis

Client:		Date Collected:	8/30/2024 12:00:00 AM
Project:		Date Received:	8/30/2024 12:00:00 AM
Client Sample ID:	PB163076BL	SDG No.:	BP091624
Lab Sample ID:	PB163076BL	Matrix:	Solid
Analytical Method:	8270E	% Moisture:	0
Sample Wt/Vol:	30.02	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
BP021937.D	1	8/30/2024 12:00:00 AM	9/16/2024 12:00:00 AM	PB163076

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
62-75-9	n-Nitrosodimethylamine	86.7	U	86.7	270	330	ug/Kg
110-86-1	Pyridine	79.8	U	79.8	130	170	ug/Kg
100-52-7	Benzaldehyde	180	U	180	270	330	ug/Kg
62-53-3	Aniline	96.8	U	96.8	130	170	ug/Kg
108-95-2	Phenol	82.8	U	82.8	130	170	ug/Kg
111-44-4	bis(2-Chloroethyl)ether	83.6	U	83.6	130	170	ug/Kg
95-57-8	2-Chlorophenol	83.4	U	83.4	130	170	ug/Kg
95-50-1	1,2-Dichlorobenzene	84.5	U	84.5	130	170	ug/Kg
541-73-1	1,3-Dichlorobenzene	85.9	U	85.9	130	170	ug/Kg
106-46-7	1,4-Dichlorobenzene	86.4	U	86.4	130	170	ug/Kg
100-51-6	Benzyl Alcohol	82.9	U	82.9	270	330	ug/Kg
95-48-7	2-Methylphenol	80.5	U	80.5	130	170	ug/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	90.8	U	90.8	130	170	ug/Kg
98-86-2	Acetophenone	86.8	U	86.8	130	170	ug/Kg
65794-96-9	3+4-Methylphenols	79.7	U	79.7	270	330	ug/Kg
621-64-7	n-Nitroso-di-n-propylamine	40.3	U	40.3	79.9	79.9	ug/Kg
67-72-1	Hexachloroethane	82.9	U	82.9	130	170	ug/Kg
98-95-3	Nitrobenzene	90.7	U	90.7	130	170	ug/Kg
78-59-1	Isophorone	84.5	U	84.5	130	170	ug/Kg
88-75-5	2-Nitrophenol	94.4	U	94.4	130	170	ug/Kg
105-67-9	2,4-Dimethylphenol	93.1	U	93.1	130	170	ug/Kg
111-91-1	bis(2-Chloroethoxy)methane	85.7	U	85.7	130	170	ug/Kg
120-83-2	2,4-Dichlorophenol	75.4	U	75.4	130	170	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	85.3	U	85.3	130	170	ug/Kg
65-85-0	Benzoic acid	240	U	240	270	330	ug/Kg
91-20-3	Naphthalene	82.5	U	82.5	130	170	ug/Kg
106-47-8	4-Chloroaniline	82.5	U	82.5	130	170	ug/Kg
87-68-3	Hexachlorobutadiene	83.2	U	83.2	130	170	ug/Kg

Report of Analysis

Client:		Date Collected:	8/30/2024 12:00:00 AM
Project:		Date Received:	8/30/2024 12:00:00 AM
Client Sample ID:	PB163076BL	SDG No.:	BP091624
Lab Sample ID:	PB163076BL	Matrix:	Solid
Analytical Method:	8270E	% Moisture:	0
Sample Wt/Vol:	30.02 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
BP021937.D	1	8/30/2024 12:00:00 AM	9/16/2024 12:00:00 AM	PB163076

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
105-60-2	Caprolactam	86.7	U	86.7	270	330	ug/Kg
59-50-7	4-Chloro-3-methylphenol	77.4	U	77.4	130	170	ug/Kg
91-57-6	2-Methylnaphthalene	82.4	U	82.4	130	170	ug/Kg
77-47-4	Hexachlorocyclopentadiene	160	U	160	270	330	ug/Kg
88-06-2	2,4,6-Trichlorophenol	71.4	U	71.4	130	170	ug/Kg
95-95-4	2,4,5-Trichlorophenol	74	U	74	130	170	ug/Kg
92-52-4	1,1-Biphenyl	87.3	U	87.3	130	170	ug/Kg
91-58-7	2-Chloronaphthalene	83.2	U	83.2	130	170	ug/Kg
88-74-4	2-Nitroaniline	94.9	U	94.9	130	170	ug/Kg
131-11-3	Dimethylphthalate	81.6	U	81.6	130	170	ug/Kg
208-96-8	Acenaphthylene	86.4	U	86.4	130	170	ug/Kg
606-20-2	2,6-Dinitrotoluene	83.1	U	83.1	130	170	ug/Kg
99-09-2	3-Nitroaniline	89.1	U	89.1	130	170	ug/Kg
83-32-9	Acenaphthene	81	U	81	130	170	ug/Kg
Total PAH	Total PAH	1323.5	U	1323.5	130	2720	ug/Kg
51-28-5	2,4-Dinitrophenol	240	U	240	270	330	ug/Kg
100-02-7	4-Nitrophenol	120	U	120	270	330	ug/Kg
132-64-9	Dibenzofuran	84.3	U	84.3	130	170	ug/Kg
25321-14-6	2,4-Dinitrotoluene/2,6-Dinitrotoluene (mixture)	169.2	U	169.2	130	340	ug/Kg
121-14-2	2,4-Dinitrotoluene	86.1	U	86.1	130	170	ug/Kg
84-66-2	Diethylphthalate	80	U	80	130	170	ug/Kg
7005-72-3	4-Chlorophenyl-phenylether	85.5	U	85.5	130	170	ug/Kg
86-73-7	Fluorene	85.4	U	85.4	130	170	ug/Kg
100-01-6	4-Nitroaniline	110	U	110	130	170	ug/Kg
534-52-1	4,6-Dinitro-2-methylphenol	120	U	120	270	330	ug/Kg
86-30-6	n-Nitrosodiphenylamine	81.5	U	81.5	130	170	ug/Kg
103-33-3	Azobenzene	79.5	U	79.5	130	170	ug/Kg
101-55-3	4-Bromophenyl-phenylether	78.8	U	78.8	130	170	ug/Kg
118-74-1	Hexachlorobenzene	84.9	U	84.9	130	170	ug/Kg
1912-24-9	Atrazine	91.3	U	91.3	130	170	ug/Kg

Report of Analysis

Client:		Date Collected:	8/30/2024 12:00:00 AM
Project:		Date Received:	8/30/2024 12:00:00 AM
Client Sample ID:	PB163076BL	SDG No.:	BP091624
Lab Sample ID:	PB163076BL	Matrix:	Solid
Analytical Method:	8270E	% Moisture:	0
Sample Wt/Vol:	30.02 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
BP021937.D	1	8/30/2024 12:00:00 AM	9/16/2024 12:00:00 AM	PB163076

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
87-86-5	Pentachlorophenol	77.2	U	77.2	270	330	ug/Kg
85-01-8	Phenanthrene	83.9	U	83.9	130	170	ug/Kg
120-12-7	Anthracene	84.3	U	84.3	130	170	ug/Kg
86-74-8	Carbazole	80.2	U	80.2	130	170	ug/Kg
84-74-2	Di-n-butylphthalate	84.2	U	84.2	130	170	ug/Kg
206-44-0	Fluoranthene	81.6	U	81.6	130	170	ug/Kg
92-87-5	Benzidine	160	U	160	270	330	ug/Kg
129-00-0	Pyrene	82.9	U	82.9	130	170	ug/Kg
85-68-7	Butylbenzylphthalate	96.7	U	96.7	130	170	ug/Kg
91-94-1	3,3-Dichlorobenzidine	98.5	U	98.5	270	330	ug/Kg
56-55-3	Benzo(a)anthracene	80.6	U	80.6	130	170	ug/Kg
218-01-9	Chrysene	79.4	U	79.4	130	170	ug/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	90.9	U	90.9	130	170	ug/Kg
117-84-0	Di-n-octyl phthalate	110	U	110	270	330	ug/Kg
205-99-2	Benzo(b)fluoranthene	81	U	81	130	170	ug/Kg
207-08-9	Benzo(k)fluoranthene	82.5	U	82.5	130	170	ug/Kg
50-32-8	Benzo(a)pyrene	92.9	U	92.9	130	170	ug/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	78	U	78	130	170	ug/Kg
53-70-3	Dibenzo(a,h)anthracene	81.1	U	81.1	130	170	ug/Kg
191-24-2	Benzo(g,h,i)perylene	80	U	80	130	170	ug/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	86.7	U	86.7	130	170	ug/Kg
123-91-1	1,4-Dioxane	110	U	110	130	170	ug/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	74.7	U	74.7	130	170	ug/Kg
90-12-0	1-Methylnaphthalene	83.1	U	83.1	170	170	ug/Kg
SURROGATES							
367-12-4	2-Fluorophenol	130.162		18-112		87	SPK: -150
13127-88-3	Phenol-d6	112.889		15-107		75	SPK: -150
4165-60-0	Nitrobenzene-d5	122.958	*	18-107		123	SPK: -100
321-60-8	2-Fluorobiphenyl	83.639		20-109		84	SPK: -100

Report of Analysis

Client:		Date Collected:	8/30/2024 12:00:00 AM
Project:		Date Received:	8/30/2024 12:00:00 AM
Client Sample ID:	PB163076BL	SDG No.:	BP091624
Lab Sample ID:	PB163076BL	Matrix:	Solid
Analytical Method:	8270E	% Moisture:	0
Sample Wt/Vol:	30.02 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
BP021937.D	1	8/30/2024 12:00:00 AM	9/16/2024 12:00:00 AM	PB163076

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
118-79-6	2,4,6-Tribromophenol	116.014		10-116		77	SPK: -150
1718-51-0	Terphenyl-d14	84.049		10-105		84	SPK: -100
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	247840	7.699				
1146-65-2	Naphthalene-d8	640886	10.475				
15067-26-2	Acenaphthene-d10	612448	14.345				
1517-22-2	Phenanthrene-d10	1343596	17.133				
1719-03-5	Chrysene-d12	1392486	21.586				
1520-96-3	Perylene-d12	1501306	24.898				
TENTATIVE IDENTIFIED COMPOUNDS							
000123-42-2	2-Pentanone, 4-hydroxy-4-methyl-	160	A			4.875	
000629-94-7	Heneicosane	87.3	J			25.045	

Hit Summary Sheet SW-846

SDG No.: BP091624

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
Client ID :	ICVBP091624							
SSTDICV040	ICVBP091624	Water	1,1-Biphenyl	30,700.000		910	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,2,4,5-Tetrachlorobenzene	29,100.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,2,4-Trichlorobenzene	39,600.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,2-Dichlorobenzene	39,200.000		880	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,3-Dichlorobenzene	39,800.000		800	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,4-Dichlorobenzene	39,400.000		840	5000	ug/L
SSTDICV040	ICVBP091624	Water	1,4-Dioxane	36,400.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	1-Methylnaphthalene	40,100.000		860	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,2-oxybis(1-Chloropropane)	38,200.000		1400	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,3,4,6-Tetrachlorophenol	41,300.000		790	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4,5-Trichlorophenol	31,000.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4,6-Trichlorophenol	31,800.000		890	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4-Dichlorophenol	40,900.000		880	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4-Dimethylphenol	42,600.000		1500	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4-Dinitrophenol	40,200.000		6400	10000	ug/L
SSTDICV040	ICVBP091624	Water	2,4-Dinitrotoluene	41,200.000		1500	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,6-Dinitrotoluene	35,800.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Chloronaphthalene	30,900.000		970	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Chlorophenol	39,000.000		710	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Methylnaphthalene	40,000.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Methylphenol	38,800.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Nitroaniline	32,100.000		1400	5000	ug/L
SSTDICV040	ICVBP091624	Water	2-Nitrophenol	44,400.000		2000	5000	ug/L
SSTDICV040	ICVBP091624	Water	3+4-Methylphenols	38,300.000		1200	10000	ug/L
SSTDICV040	ICVBP091624	Water	3,3-Dichlorobenzidine	41,700.000		1300	10000	ug/L
SSTDICV040	ICVBP091624	Water	3-Nitroaniline	40,700.000		1400	5000	ug/L
SSTDICV040	ICVBP091624	Water	4,6-Dinitro-2-methylphenol	40,500.000		3100	10000	ug/L
SSTDICV040	ICVBP091624	Water	4-Bromophenyl-phenylether	39,800.000		950	5000	ug/L
SSTDICV040	ICVBP091624	Water	4-Chloro-3-methylphenol	40,700.000		840	5000	ug/L
SSTDICV040	ICVBP091624	Water	4-Chloroaniline	38,800.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	4-Chlorophenyl-phenylether	39,700.000		980	5000	ug/L
SSTDICV040	ICVBP091624	Water	4-Nitroaniline	40,500.000		2000	5000	ug/L
SSTDICV040	ICVBP091624	Water	4-Nitrophenol	41,100.000		2000	10000	ug/L
SSTDICV040	ICVBP091624	Water	Acenaphthene	38,600.000		810	5000	ug/L
SSTDICV040	ICVBP091624	Water	Acenaphthylene	40,000.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	Acetophenone	40,200.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Aniline	38,300.000		1600	5000	ug/L

Hit Summary Sheet SW-846

SDG No.: BP091624

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
SSTDICV040	ICVBP091624	Water	Anthracene	39,800.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Atrazine	38,600.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	Azobenzene	39,300.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzaldehyde	43,400.000		4000	10000	ug/L
SSTDICV040	ICVBP091624	Water	Benzidine	34,400.000		4100	10000	ug/L
SSTDICV040	ICVBP091624	Water	Benzo(a)anthracene	39,500.000		940	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzo(a)pyrene	39,700.000		1700	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzo(b)fluoranthene	40,400.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzo(g,h,i)perylene	39,400.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzo(k)fluoranthene	41,700.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Benzoic acid	42,500.000		5500	10000	ug/L
SSTDICV040	ICVBP091624	Water	Benzyl Alcohol	38,800.000		1600	10000	ug/L
SSTDICV040	ICVBP091624	Water	bis(2-Chloroethoxy)methane	39,500.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	bis(2-Chloroethyl)ether	38,700.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Bis(2-ethylhexyl)phthalate	39,600.000		1900	5000	ug/L
SSTDICV040	ICVBP091624	Water	Butylbenzylphthalate	39,200.000		2100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Caprolactam	39,800.000		1700	10000	ug/L
SSTDICV040	ICVBP091624	Water	Carbazole	40,300.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Chrysene	38,900.000		860	5000	ug/L
SSTDICV040	ICVBP091624	Water	Di-n-butylphthalate	34,900.000		1500	5000	ug/L
SSTDICV040	ICVBP091624	Water	Di-n-octyl phthalate	32,200.000		2500	10000	ug/L
SSTDICV040	ICVBP091624	Water	Dibenzo(a,h)anthracene	38,900.000		1200	5000	ug/L
SSTDICV040	ICVBP091624	Water	Dibenzofuran	40,900.000		930	5000	ug/L
SSTDICV040	ICVBP091624	Water	Diethylphthalate	37,900.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	Dimethylphthalate	33,300.000		930	5000	ug/L
SSTDICV040	ICVBP091624	Water	Fluoranthene	39,500.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	Fluorene	40,200.000		960	5000	ug/L
SSTDICV040	ICVBP091624	Water	Hexachlorobenzene	39,400.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Hexachlorobutadiene	40,700.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	Hexachlorocyclopentadiene	30,400.000		5000	10000	ug/L
SSTDICV040	ICVBP091624	Water	Hexachloroethane	40,400.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	Indeno(1,2,3-cd)pyrene	39,300.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	Isophorone	42,600.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	n-Nitroso-di-n-propylamine	37,900.000		1500	2500	ug/L
SSTDICV040	ICVBP091624	Water	n-Nitrosodimethylamine	36,000.000		1000	10000	ug/L
SSTDICV040	ICVBP091624	Water	n-Nitrosodiphenylamine	40,600.000		890	5000	ug/L
SSTDICV040	ICVBP091624	Water	Naphthalene	39,200.000		1000	5000	ug/L
SSTDICV040	ICVBP091624	Water	Nitrobenzene	41,000.000		1300	5000	ug/L
SSTDICV040	ICVBP091624	Water	Pentachlorophenol	40,300.000		1900	10000	ug/L
SSTDICV040	ICVBP091624	Water	Phenanthrene	39,700.000		890	5000	ug/L

Hit Summary Sheet
SW-846

SDG No.: BP091624

Client:

Sample ID	Client ID		Parameter	Concentration	C	MDL	RDL	Units
SSTDICV040	ICVBP091624	Water	Phenol	38,900.000		930	5000	ug/L
SSTDICV040	ICVBP091624	Water	Pyrene	40,900.000		1100	5000	ug/L
SSTDICV040	ICVBP091624	Water	Pyridine	40,200.000		1600	5000	ug/L
SSTDICV040	ICVBP091624	Water	2,4-Dinitrotoluene/2,6-Dinitrotolu	77,000.000		2700	10000	ug/L
SSTDICV040	ICVBP091624	Water	Total PAH	635,700.000		17360	80000	ug/L
Total Svoc :				3,812,900.00				
Total Concentration:				3,812,900.00				



6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: _____
Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG No.: BP091624
Instrument ID: _____ Calibration Date(s): 9/16/2024 : _____
Calibration Time(s): 11:10 _____

LAB FILE ID:		RRF2.5 = BP021928.D			RRF005 = BP021929.D		RRF010 = BP021930.D	
		RRF020 = BP021931.D			RRF040 = BP021932.D		RRF050 = BP021933.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
n-Nitrosodimethylamine		0.717	0.716	0.637	0.681	0.442	0.669	19.1
Pyridine		1.148	1.168	1.171	1.248	0.849	1.117	13.8
2-Fluorophenol		1.061	1.041	1.040	1.058	1.043	1.108	14.6
Benzaldehyde		0.768	0.834	0.841	0.771	0.626	0.724	15.3
Aniline		1.203	1.199	1.221	1.246	1.259	1.227	2.7
Phenol-d6		1.363	1.344	1.362	1.390	1.190	1.358	6.0
Phenol		1.378	1.333	1.369	1.385	1.178	1.352	6.0
bis(2-Chloroethyl)ether		1.083	1.053	1.097	1.106	1.105	1.084	1.9
2-Chlorophenol		1.186	1.169	1.147	1.161	1.117	1.155	2.0
1,2-Dichlorobenzene		1.458	1.454	1.396	1.391	1.334	1.392	3.6
1,3-Dichlorobenzene		1.520	1.471	1.439	1.450	1.387	1.437	3.7
1,4-Dichlorobenzene		1.513	1.494	1.476	1.455	1.408	1.455	3.0
Benzyl Alcohol		0.990	1.014	1.038	1.026	1.003	1.042	4.9
2-Methylphenol		0.919	0.915	0.939	0.962	0.934	0.944	2.5
2,2-oxybis(1-Chloropropane)		1.137	1.105	1.206	1.268	1.357	1.200	7.3
Acetophenone		0.499	0.504	0.498	0.359	0.468	0.470	10.9
3+4-Methylphenols		1.225	1.277	1.321	1.357	1.326	1.324	4.3
n-Nitroso-di-n-propylamine	0.813	0.917	0.896	0.935	0.944	0.904	0.913	5.0
Nitrobenzene-d5		0.385	0.379	0.380	0.281	0.362	0.365	10.5
Hexachloroethane		0.544	0.552	0.530	0.557	0.500	0.535	3.8
Nitrobenzene		0.408	0.390	0.390	0.292	0.375	0.377	10.4
Isophorone		0.602	0.610	0.638	0.486	0.441	0.583	14.6
2-Nitrophenol		0.137	0.144	0.155	0.122	0.139	0.149	12.7
2,4-Dimethylphenol		0.185	0.189	0.197	0.153	0.168	0.185	9.8
bis(2-Chloroethoxy)methane		0.373	0.372	0.398	0.397	0.396	0.386	3.4
2,4-Dichlorophenol		0.292	0.303	0.321	0.321	0.320	0.316	4.4
1,2,4-Trichlorobenzene		0.412	0.388	0.385	0.384	0.377	0.385	3.7
Benzoic acid		0.171	0.188	0.192	0.165	0.215	0.202	15.4
Naphthalene		1.046	1.011	1.019	1.002	0.964	1.000	3.2
4-Chloroaniline		0.354	0.367	0.374	0.372	0.367	0.368	2.4
Hexachlorobutadiene		0.296	0.283	0.266	0.253	0.243	0.266	7.1
Caprolactam		0.083	0.090	0.095	0.101	0.101	0.096	7.5
4-Chloro-3-methylphenol		0.320	0.339	0.352	0.338	0.336	0.343	4.3
2-Methylnaphthalene		0.751	0.742	0.739	0.718	0.703	0.727	2.9
Hexachlorocyclopentadiene		0.205	0.211	0.206	0.199	0.203	0.209	4.5
2,4,6-Trichlorophenol		0.363	0.350	0.375	0.390	0.394	0.384	5.7

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1



6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: _____
Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG No.: BP091624
Instrument ID: _____ Calibration Date(s): 9/16/2024 : _____
Calibration Time(s): 11:10 _____

LAB FILE ID:		RRF2.5 = BP021928.D			RRF005 = BP021929.D		RRF010 = BP021930.D	
		RRF020 = BP021931.D			RRF040 = BP021932.D		RRF050 = BP021933.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
2-Fluorobiphenyl		1.429	1.313	1.323	1.268	1.248	1.312	4.7
2,4,5-Trichlorophenol		0.432	0.416	0.433	0.439	0.449	0.440	3.5
1,1-Biphenyl		1.469	1.353	1.386	1.339	1.332	1.363	3.9
2-Chloronaphthalene		1.159	1.090	1.091	1.081	1.071	1.094	2.9
2-Nitroaniline		0.253	0.276	0.301	0.309	0.305	0.301	9.3
Dimethylphthalate		1.539	1.448	1.451	1.392	1.381	1.432	4.0
Acenaphthylene		1.574	1.538	1.593	1.566	1.537	1.567	1.8
2,6-Dinitrotoluene		0.299	0.302	0.318	0.319	0.321	0.315	3.5
3-Nitroaniline		0.257	0.268	0.283	0.302	0.303	0.289	7.0
Acenaphthene		1.054	1.028	1.031	1.002	0.993	1.021	2.9
2,4-Dinitrophenol		0.129	0.146	0.157	0.164	0.170	0.164	13.7
4-Nitrophenol		0.213	0.230	0.258	0.268	0.265	0.252	8.6
Dibenzofuran		1.818	1.739	1.735	1.674	1.135	1.631	13.9
2,4-Dinitrotoluene		0.414	0.426	0.452	0.444	0.327	0.426	11.1
Diethylphthalate		1.587	1.554	1.503	1.393	0.978	1.414	14.5
4-Chlorophenyl-phenylether		0.820	0.803	0.747	0.725	0.472	0.716	16.1
Fluorene		1.495	1.470	1.421	1.365	0.883	1.335	15.6
4-Nitroaniline		0.295	0.320	0.320	0.322	0.218	0.305	13.2
4,6-Dinitro-2-methylphenol		0.095	0.097	0.102	0.106	0.103	0.106	9.1
n-Nitrosodiphenylamine		0.523	0.525	0.533	0.506	0.469	0.512	4.2
Azobenzene		1.294	1.350	1.328	1.273	0.855	1.234	13.9
2,4,6-Tribromophenol		0.279	0.287	0.269	0.254	0.191	0.263	12.8
4-Bromophenyl-phenylether		0.204	0.207	0.209	0.199	0.198	0.204	2.2
Hexachlorobenzene		0.252	0.247	0.243	0.232	0.236	0.243	2.8
Atrazine		0.199	0.200	0.198	0.185	0.182	0.187	6.4
Pentachlorophenol		0.149	0.152	0.164	0.160	0.169	0.163	6.4
Phenanthrene		1.046	0.984	1.004	0.974	0.944	0.983	3.5
Anthracene		0.958	0.952	0.974	0.941	0.924	0.944	2.2
Carbazole		0.959	0.939	0.984	0.939	0.926	0.936	3.3
Di-n-butylphthalate		1.052	1.070	1.119	1.073	1.140	1.092	3.0
Fluoranthene		1.318	1.291	1.322	1.290	1.289	1.283	2.8
Benzidine		0.292	0.378	0.255	0.228	0.302	0.291	19.6
Pyrene		1.207	1.267	1.258	1.312	1.266	1.278	3.2
Terphenyl-d14		1.005	1.059	0.986	0.983	0.983	1.013	3.1
Butylbenzylphthalate		0.421	0.454	0.480	0.482	0.506	0.483	7.6
3,3-Dichlorobenzidine		0.406	0.416	0.442	0.441	0.447	0.429	3.9

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1



6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: _____
Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG No.: BP091624
Instrument ID: _____ Calibration Date(s): 9/16/2024 : _____
Calibration Time(s): 11:10 _____

LAB FILE ID:		RRF2.5 = BP021928.D			RRF005 = BP021929.D		RRF010 = BP021930.D	
		RRF020 = BP021931.D			RRF040 = BP021932.D		RRF050 = BP021933.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
Benzo(a)anthracene		1.291	1.293	1.315	1.276	1.266	1.288	1.3
Chrysene		1.292	1.231	1.247	1.221	1.195	1.232	2.6
Bis(2-ethylhexyl)phthalate		0.622	0.672	0.719	0.711	0.744	0.705	6.2
Di-n-octyl phthalate		0.973	1.027	1.169	1.172	1.277	1.163	10.4
Benzo(b)fluoranthene		1.138	1.126	0.935	1.179	1.201	1.143	8.6
Benzo(k)fluoranthene		1.138	1.105	0.865	1.138	1.119	1.085	9.1
Benzo(a)pyrene		0.988	0.973	1.020	1.017	1.032	1.016	2.6
Indeno(1,2,3-cd)pyrene		1.348	1.290	1.265	1.258	1.236	1.284	2.8
Dibenzo(a,h)anthracene		1.130	1.081	1.060	1.031	1.017	1.063	3.5
Benzo(g,h,i)perylene		1.135	1.121	1.090	1.090	1.077	1.106	2.0
1,2,4,5-Tetrachlorobenzene		0.631	0.578	0.600	0.590	0.586	0.600	3.1
1,4-Dioxane		0.477	0.450	0.429	0.615	0.431	0.481	16.2
2,3,4,6-Tetrachlorophenol		0.395	0.388	0.390	0.391	0.259	0.375	13.7
1-Methylnaphthalene		0.744	0.737	0.731	0.715	0.701	0.721	2.7

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: BP091624 SAS No.: BP091624 SDG NO.: BP091624

EPA Sample No.: SSTDICCC040 Date Analyzed: Sep 16 2024 12:00AM

Lab File ID: BP021932.D Time Analyzed: 16:35

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

	IS1 (DCB) AREA #	RT #	IS2 (NPT) AREA #	RT #	IS3 (ANT) AREA #	RT #
12 HOUR STD	118788	7.746	626099	10.51	463586	14.35
UPPER LIMIT	237576	8.246	1252198	11.01	927172	14.851
LOWER LIMIT	59394	7.246	313049.5	10.01	231793	13.851
EPA SAMPLE NO.						
01 ICVBP091624	120317	7.75	444569	10.51	465447	14.36
02 PB163076BL	247840 *	7.70	640886	10.48	612448	14.35

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.: BP091624 SAS No.: BP091624 SDG NO.: BP091624

EPA Sample No.: SSTDICCC040 Date Analyzed: Sep 16 2024 12:00AM

Lab File ID: BP021932.D Time Analyzed: 16:35

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

	IS1 (DCB) AREA #	RT #	IS2 (NPT) AREA #	RT #	IS3 (ANT) AREA #	RT #
12 HOUR STD	118788	7.746	626099	10.51	463586	14.35
UPPER LIMIT	237576	8.246	1252198	11.01	927172	14.851
LOWER LIMIT	59394	7.246	313049.5	10.01	231793	13.851
EPA SAMPLE NO.						
01 ICVBP091624	120317	7.75	444569	10.51	465447	14.36
02 PB163076BL	247840 *	7.70	640886	10.48	612448	14.35

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.: BP091624 SAS No.: BP091624 SDG NO.: BP091624

EPA Sample No.: SSTDICCC040 Date Analyzed: Sep 16 2024 12:00AM

Lab File ID: BP021932.D Time Analyzed: 16:35

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

	IS4 (PHN) AREA #	RT #	IS5 (CRY) AREA #	RT #	IS6 (PRY) AREA #	RT #
12 HOUR STD	1.07216e+04	17.139	1.11829e+	21.575	1.2651e+004	24.868
UPPER LIMIT	2144320	17.639	2236580	22.075	2530200	25.368
LOWER LIMIT	536080	16.639	559145	21.075	632550	24.368
EPA SAMPLE NO.						
01 ICVBP091624	1.02561	17.15	1.04969e+	21.59	1.18952e+	24.89
02 PB163076BL	1.3436e	17.13	1.39249e+	21.59	1.50131e+	24.90

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.: BP091624 SAS No.: BP091624 SDG NO.: BP091624

EPA Sample No.: SSTDICCC040 Date Analyzed: Sep 16 2024 12:00AM

Lab File ID: BP021932.D Time Analyzed: 16:35

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

	IS4 (PHN) AREA #	RT #	IS5 (CRY) AREA #	RT #	IS6 (PRY) AREA #	RT #
12 HOUR STD	1.07216e+04	17.139	1.11829e+	21.575	1.2651e+004	24.868
UPPER LIMIT	2144320	17.639	2236580	22.075	2530200	25.368
LOWER LIMIT	536080	16.639	559145	21.075	632550	24.368
EPA SAMPLE NO.						
01 ICVBP091624	1.02561	17.15	1.04969e+	21.59	1.18952e+	24.89
02 PB163076BL	1.3436e	17.13	1.39249e+	21.59	1.50131e+	24.90

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.

ICVBP091624

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BP091624

SAS No.: BP091624 SDG NO.: BP091624

Lab File ID: BP021936.D

Lab Sample ID: SSTDICV040

Instrument ID: BNA_P

Date Extracted: _____

Matrix: (soil/water) Water

Date Analyzed: 09/16/2024

Level: (low/med) LOW

Time Analyzed: 16:35

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
ICVBP091624	SSTDICV040	BP021936.D	09/16/2024

COMMENTS: _____

4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

PB163076BL

Lab Name: CHEMTECH

Contract: _____

Lab Code: CHEM Case No.: BP091624

SAS No.: BP091624 SDG NO.: BP091624

Lab File ID: BP021937.D

Lab Sample ID: PB163076BL

Instrument ID: BNA_P

Date Extracted: 08/30/2024

Matrix: (soil/water) Solid

Date Analyzed: 09/16/2024

Level: (low/med) LOW

Time Analyzed: 17:35

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.: BP091624

Client: _____

Analytical Method: 8270E

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
SSTDICV040	ICVBP091624	BP021936.D	2-Fluorophenol	150	75,094.00	50063	*	10	139
			Phenol-d6	150	78,812.00	52541	*	10	134
			Nitrobenzene-d5	100	82,446.00	82446	*	49	133
			2-Fluorobiphenyl	100	60,600.00	60600	*	52	132
			2,4,6-Tribromophenol	150	76,181.00	50787	*	44	137
			Terphenyl-d14	100	78,206.00	78206	*	48	125

Surrogate Summary

SW-846

SDG No.: BP091624

Client: _____

Analytical Method: 8270E

Lab Sample ID	Client ID	Datafile	Parameter	Spike (PPM)	Result (PPM)	Recovery (%)	Qual	Limits (%)	
								Low	High
PB163076BL	PB163076BL	BP021937.D	2-Fluorophenol	150	130.16	87	*	18	112
			Phenol-d6	150	112.89	75		15	107
			Nitrobenzene-d5	100	122.96	123		18	107
			2-Fluorobiphenyl	100	83.64	84		20	109
			2,4,6-Tribromophenol	150	116.01	77		10	116
			Terphenyl-d14	100	84.05	84		10	105

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: CHEMTECH

Contract: BP091624

Lab Code: CHEM

SAS No.: BP091624 SDG NO.: BP091624

Lab File ID: BP021927.D

DFTPP Injection Date: 09/16/2024

Instrument ID: BNA_P

DFTPP Injection Time: 10:30

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	34.8
68	Less than 2.0% of mass 69	0.7 (1.8) 1
69	Mass 69 relative abundance	37.3
70	Less than 2.0% of mass 69	0.2 (0.5) 1
127	10.0 - 80.0% of mass 198	43.4
197	Less than 2.0% of mass 198	0.7
198	Base Peak, 100% relative abundance	100
199	5.0 to 9.0% of mass 198	6.8
275	10.0 - 60.0% of mass 198	30.1
365	Greater than 1% of mass 198	6
441	Present, but less than mass 443	12.9
442	Greater than 50% of mass 198	83.5
443	15.0 - 24.0% of mass 442	15.9 (19.1) 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
SSTDICC2.5	SSTDICC2.5	BP021928.D	09/16/2024	11:10
SSTDICC005	SSTDICC005	BP021929.D	09/16/2024	11:51
SSTDICC010	SSTDICC010	BP021930.D	09/16/2024	12:32
SSTDICC020	SSTDICC020	BP021931.D	09/16/2024	13:12
SSTDICCC040	SSTDICCC040	BP021932.D	09/16/2024	13:53
SSTDICC050	SSTDICC050	BP021933.D	09/16/2024	14:33
SSTDICC060	SSTDICC060	BP021934.D	09/16/2024	15:14
SSTDICC080	SSTDICC080	BP021935.D	09/16/2024	15:55
ICVBP091624	SSTDICV040	BP021936.D	09/16/2024	16:35
PB163076BL	PB163076BL	BP021937.D	09/16/2024	17:35