

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090924\
 Data File : BG062688.D
 Acq On : 9 Sep 2024 14:21
 Operator : JU/RC
 Sample : SST00552
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SST005438

Quant Time: Sep 10 01:00:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 00:41:25 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.908	152	66293	20.000	ng/u1	0.00
20) Naphthalene-d8	10.699	136	281739	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.536	164	224986	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.279	188	600173	20.000	ng/u1	0.00
79) Chrysene-d12	21.522	240	640624	20.000	ng/u1	0.00
88) Perylene-d12	24.589	264	676984	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.366	96	2469	1.676	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	0.000	99	0d	0.000	ng/u1	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/u1	
11) 2-Chlorophenol-d4	7.438	132	19256	4.548	ng/u1	0.00
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/u1	
21) Nitrobenzene-d5	9.054	128	8564	4.253	ng/u1	0.00
24) 2-Nitrophenol-d4	9.776	143	8858	4.022	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.323	165	21871	4.560	ng/u1	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/u1	
46) Dimethylphthalate-d6	13.936	166	83733	4.833	ng/u1	0.00
49) Acenaphthylene-d8	14.224	160	89721	4.663	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.523	176	74769	4.789	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/u1	
73) Anthracene-d10	17.373	188	132706	4.685	ng/u1	0.00
81) Pyrene-d10	19.665	212	172284	4.779	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.371	264	163278	4.568	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.402	88	2954	1.797	ng/uL#	82
12) 2-Chlorophenol	7.467	128	19357	4.471	ng/u1	95
17) N-Nitroso-di-n-propyla...	8.707	70	17455	4.235	ng/u1	95
19) Hexachloroethane	8.989	117	7826	4.490	ng/u1	96
22) Nitrobenzene	9.101	77	23645	4.341	ng/u1#	94
23) Isophorone	9.624	82	50429	4.528	ng/u1	100
25) 2-Nitrophenol	9.812	139	9959	4.103	ng/u1	95
26) 2,4-Dimethylphenol	9.870	107	23343	4.501	ng/u1	95
27) Bis(2-Chloroethoxy)met...	10.105	93	28926	4.647	ng/u1#	96
29) 2,4-Dichlorophenol	10.346	162	20589	4.359	ng/u1#	86
30) Naphthalene	10.746	128	71568	4.738	ng/u1	97
33) Hexachlorobutadiene	11.040	225	18737	4.762	ng/u1	88
35) 4-Chloro-3-methylphenol	11.980	107	23211	4.505	ng/u1	97
36) 2-Methylnaphthalene	12.356	142	52306	4.692	ng/u1	99
37) 1-Methylnaphthalene	12.573	142	54807	4.799	ng/u1	95
39) 1,2,4,5-Tetrachloroben...	12.726	216	35912	4.643	ng/u1	96
41) 2,4,6-Trichlorophenol	12.967	196	21339	4.555	ng/u1	98
42) 2,4,5-Trichlorophenol	13.037	196	21856	4.345	ng/u1	96
43) 1,1'-Biphenyl	13.366	154	75888	4.797	ng/u1	97
44) 2-Chloronaphthalene	13.407	162	60776	4.724	ng/u1	99
45) 2-Nitroaniline	13.607	65	14421	4.085	ng/u1	96
47) Dimethylphthalate	13.983	163	84606	4.742	ng/u1	99
48) 2,6-Dinitrotoluene	14.101	165	12280	3.907	ng/u1#	85
50) Acenaphthylene	14.254	152	96452	4.748	ng/u1	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090924\
 Data File : BG062688.D
 Acq On : 9 Sep 2024 14:21
 Operator : JU/RC
 Sample : SST00552
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SST005438

Quant Time: Sep 10 01:00:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 00:41:25 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Acenaphthene	14.594	153	68994	4.795	ng/ul	96
56) Dibenzofuran	14.929	168	102863	4.812	ng/ul	99
57) 2,4-Dinitrotoluene	14.894	165	20259	4.122	ng/ul#	94
58) 2,3,4,6-Tetrachlorophenol	15.158	232	22654	4.444	ng/ul	95
59) Diethylphthalate	15.352	149	83026	4.717	ng/ul	95
61) Fluorene	15.581	166	83604	4.979	ng/ul	93
62) 4-Chlorophenyl-phenyle...	15.576	204	48199	4.893	ng/ul	96
67) N-Nitrosodiphenylamine	15.787	169	72428	4.705	ng/ul	95
68) 4-Bromophenyl-phenylether	16.469	248	32456	4.699	ng/ul	96
69) Hexachlorobenzene	16.586	284	37928	4.810	ng/ul	92
72) Phenanthrene	17.321	178	150687	4.828	ng/ul	99
74) Anthracene	17.409	178	151950	4.776	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.331	216	37176	4.536	ng/ul	89
76) Pentachlorobenzene	14.853	250	44178	4.864	ng/ul	94
78) Di-n-butylphthalate	18.243	149	143390	4.595	ng/ul	98
80) Fluoranthene	19.336	202	191224	4.789	ng/ul	100
82) Pyrene	19.694	202	211179	4.914	ng/ul	99
83) Butylbenzylphthalate	20.587	149	54221	4.250	ng/ul	95
85) Benzo(a)anthracene	21.504	228	215109	4.828	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.427	149	84138	4.479	ng/ul	95
87) Chrysene	21.569	228	204996	4.850	ng/ul	99
90) Benzo(b)fluoranthene	23.613	252	198372	4.746	ng/ul	97
91) Benzo(k)fluoranthene	23.678	252	200678	4.689	ng/ul	97
93) Benzo(a)pyrene	24.436	252	183605	4.657	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	28.026	276	221558	4.561	ng/ul	100
95) Dibenzo(a,h)anthracene	28.090	278	175449	4.497	ng/ul#	95
96) Benzo(g,h,i)perylene	29.107	276	178620	4.761	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quant Time: Sep 10 01:00:59 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 00:41:25 2024
Response via : Initial Calibration

