

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061617\
 Data File : BG027520.D
 Acq On : 17 Jun 2017 23:47
 Operator : SJ/MA
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD02088

Quant Time: Jun 19 01:55:34 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 19 01:52:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.52	152	40373	20.00	ng/ul	0.00
18) Naphthalene-d8	11.34	136	207020	20.00	ng/ul	0.00
35) Acenaphthene-d10	15.10	164	135245	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.85	188	318750	20.00	ng/ul	0.00
75) Chrysene-d12	22.23	240	341574	20.00	ng/ul	0.00
83) Perylene-d12	25.89	264	309249	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	4.08	96	7412	7.34	ng/uL	0.00
5) Phenol-d5	7.65	99	90231	20.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.83	67	62476	20.94	ng/ul	0.00
9) 2-Chlorophenol-d4	8.05	132	64106	22.16	ng/ul	0.00
13) 4-Methylphenol-d8	9.19	113	72111	21.07	ng/ul	0.00
19) Nitrobenzene-d5	9.68	128	31906	21.22	ng/ul	0.00
22) 2-Nitrophenol-d4	10.40	143	35176	22.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.94	165	65209	21.60	ng/ul	0.00
29) 4-Chloroaniline-d4	11.46	131	91282	23.89	ng/ul	0.00
43) Dimethylphthalate-d6	14.49	166	220282	19.45	ng/ul	0.00
46) Acenaphthylene-d8	14.80	160	269259	20.65	ng/ul	0.00
51) 4-Nitrophenol-d4	15.28	143	40112	18.11	ng/ul	0.00
57) Fluorene-d10	16.09	176	207925	19.57	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.20	200	37626	19.01	ng/ul	0.00
70) Anthracene-d10	17.95	188	301269	20.27	ng/ul	0.00
76) Pyrene-d10	20.22	212	328021	19.66	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.63	264	295815	21.24	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	4.11	88	8456	7.236	ng/uL 93
4) Benzaldehyde	7.66	77	64835	25.328	ng/ul 91
6) Phenol	7.68	94	96123	20.951	ng/ul# 88
8) Bis(2-Chloroethyl)ether	7.92	93	67473	19.998	ng/ul 97
10) 2-Chlorophenol	8.08	128	63112	21.674	ng/ul 95
11) 2-Methylphenol	8.94	108	67458	20.339	ng/ul 99
12) 2,2'-oxybis(1-Chloropropan	9.04	45	121539	21.640	ng/ul 98
14) Acetophenone	9.34	105	112440	21.254	ng/ul# 89
15) N-Nitroso-di-n-propylamine	9.31	70	66828	22.359	ng/ul# 86
16) 4-Methylphenol	9.26	108	75777	20.805	ng/ul 99
17) Hexachloroethane	9.61	117	23749	21.273	ng/ul# 81
20) Nitrobenzene	9.72	77	92121	21.749	ng/ul# 94
21) Isophorone	10.24	82	178045	21.315	ng/ul# 95
23) 2-Nitrophenol	10.43	139	37369	21.299	ng/ul# 81
24) 2,4-Dimethylphenol	10.47	107	89665	21.455	ng/ul 95
25) Bis(2-Chloroethoxy)methane	10.71	93	103593	20.160	ng/ul 94
27) 2,4-Dichlorophenol	10.97	162	65385	21.431	ng/ul 98
28) Naphthalene	11.39	128	214043	20.119	ng/ul 98
30) 4-Chloroaniline	11.49	127	88696	23.523	ng/ul 94
31) Hexachlorobutadiene	11.66	225	39700	21.154	ng/ul 91
32) Caprolactam	12.23	113	26566	19.743	ng/ul 94
33) 4-Chloro-3-methylphenol	12.57	107	88955	22.137	ng/ul 98
34) 2-Methylnaphthalene	12.95	142	162267	20.568	ng/ul 98

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061617\
 Data File : BG027520.D
 Acq On : 17 Jun 2017 23:47
 Operator : SJ/MA
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD02088

Quant Time: Jun 19 01:55:34 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 19 01:52:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	13.31	216	84471	21.099	ng/ul	95
37) Hexachlorocyclopentadiene	13.29	237	37766	16.843	ng/ul	98
38) 2,4,6-Trichlorophenol	13.54	196	57277	21.620	ng/ul	94
39) 2,4,5-Trichlorophenol	13.61	196	59589	21.601	ng/ul	96
40) 1,1'-Biphenyl	13.94	154	212350	19.847	ng/ul#	96
41) 2-Chloronaphthalene	13.99	162	164757	20.418	ng/ul	94
42) 2-Nitroaniline	14.18	65	69065	22.749	ng/ul	99
44) Dimethylphthalate	14.53	163	218226	19.430	ng/ul	98
45) 2,6-Dinitrotoluene	14.66	165	44016	21.051	ng/ul#	94
47) Acenaphthylene	14.83	152	262310	20.069	ng/ul	99
48) 3-Nitroaniline	15.00	138	45689	20.136	ng/ul#	89
49) Acenaphthene	15.17	153	182808	20.054	ng/ul	96
50) 2,4-Dinitrophenol	15.20	184	25349	17.499	ng/ul	96
52) 4-Nitrophenol	15.29	109	35170	19.962	ng/ul#	56
53) Dibenzofuran	15.50	168	262333	19.744	ng/ul	98
54) 2,4-Dinitrotoluene	15.45	165	65009	20.307	ng/ul#	90
55) 2,3,4,6-Tetrachlorophenol	15.72	232	56595	21.094	ng/ul#	94
56) Diethylphthalate	15.89	149	227493	19.166	ng/ul	99
58) Fluorene	16.15	166	215945	19.585	ng/ul	97
59) 4-Chlorophenyl-phenylether	16.13	204	110867	20.005	ng/ul	90
60) 4-Nitroaniline	16.16	138	49150	17.685	ng/ul#	70
63) 4,6-Dinitro-2-methylphenol	16.22	198	39516	18.922	ng/ul#	98
64) N-Nitrosodiphenylamine	16.34	169	190770	20.806	ng/ul	96
65) 4-Bromophenyl-phenylether	17.03	248	69375	21.488	ng/ul	90
66) Hexachlorobenzene	17.16	284	72706	21.378	ng/ul	96
67) Atrazine	17.28	200	70656	20.119	ng/ul	98
68) Pentachlorophenol	17.50	266	41872	18.552	ng/ul	95
69) Phenanthrene	17.90	178	339443	19.817	ng/ul	99
71) Anthracene	17.99	178	353274	20.283	ng/ul	99
72) Carbazole	18.25	167	294961	19.690	ng/ul	99
73) Di-n-butylphthalate	18.78	149	376766	20.339	ng/ul#	97
74) Fluoranthene	19.89	202	387011	20.081	ng/ul#	87
77) Pyrene	20.25	202	395086	19.568	ng/ul#	85
78) Butylbenzylphthalate	21.13	149	169142	22.229	ng/ul	95
79) 3,3'-Dichlorobenzidine	22.10	252	137302	21.837	ng/ul#	95
80) Benzo(a)anthracene	22.21	228	392031	20.610	ng/ul	99
81) Bis(2-ethylhexyl)phthalate	22.06	149	233930	21.406	ng/ul	99
82) Chrysene	22.28	228	358351	20.860	ng/ul	99
84) Di-n-octyl phthalate	23.42	149	405948	20.950	ng/ul	100
85) Benzo(b)fluoranthene	24.71	252	379900	20.626	ng/ul#	95
86) Benzo(k)fluoranthene	24.79	252	365035	20.619	ng/ul#	95
88) Benzo(a)pyrene	25.72	252	367830	20.800	ng/ul#	96
89) Indeno(1,2,3-cd)pyrene	30.08	276	414110	20.513	ng/ul#	90
90) Dibenzo(a,h)anthracene	30.16	278	359104	20.718	ng/ul#	93
91) Benzo(g,h,i)perylene	31.42	276	306930	18.570	ng/ul#	92

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

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061617\
Data File : BG027520.D
Acq On : 17 Jun 2017 23:47
Operator : SJ/MA
Sample : SSTDCCC020EC
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
Client Sampled :
SSTD02088

Quant Time: Jun 19 01:55:34 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 19 01:52:26 2017
Response via : Initial Calibration

