

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082218\
 Data File : BG036340.D
 Acq On : 22 Aug 2018 13:22
 Operator : JU/SJ
 Sample : MDL-W-07
 Misc : 1PPM/4PPM
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-W-07

Quant Time: Aug 22 13:59:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG082118MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 22 10:52:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	52835	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	224166	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.70	164	145058	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.44	188	357674	20.00	ng/ul	0.00
78) Chrysene-d12	21.71	240	384809	20.00	ng/ul	0.00
86) Perylene-d12	24.94	264	374279	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.51	96	8355	5.89	ng/uL	0.00
5) Phenol-d5	7.22	99	152096	30.94	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	107068	31.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.60	132	109613	30.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.77	113	121283	32.32	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	52567	38.44	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	51869	37.70	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.49	165	113849	30.91	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	160389	39.37	ng/ul	0.00
44) Dimethylphthalate-d6	14.10	166	395715	32.62	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	473097	32.35	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	60288	33.60	ng/ul	0.00
58) Fluorene-d10	15.69	176	328626	31.16	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	39717	30.66	ng/ul	0.00
71) Anthracene-d10	17.44	188	357674	21.43	ng/ul	-0.10
79) Pyrene-d10	19.82	212	626641	31.78	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.72	264	597567	29.42	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.55	88	2518	1.683 ng/uL#	27
4) Benzaldehyde	7.21	77	7834	2.378 ng/ul	99
6) Phenol	7.24	94	18505	3.767 ng/ul#	86
8) Bis(2-Chloroethyl)ether	7.49	93	15027	4.119 ng/ul#	86
10) 2-Chlorophenol	7.63	128	13209	3.721 ng/ul#	79
11) 2-Methylphenol	8.51	108	13770	3.679 ng/ul	89
12) 2,2'-oxybis(1-Chloropropan	8.60	45	30795	3.834 ng/ul#	88
14) Acetophenone	8.90	105	23267	4.045 ng/ul#	90
15) N-Nitroso-di-n-propylamine	8.88	70	13078	3.774 ng/ul#	90
16) 4-Methylphenol	8.82	108	13948	3.562 ng/ul	96
17) Hexachloroethane	9.17	117	5725	4.048 ng/ul#	77
20) Nitrobenzene	9.28	77	17278	4.453 ng/ul#	96
21) Isophorone	9.80	82	35453	3.764 ng/ul	98
23) 2-Nitrophenol	9.98	139	6659	4.573 ng/ul	100
24) 2,4-Dimethylphenol	10.04	107	17579	3.952 ng/ul	89
25) Bis(2-Chloroethoxy)methane	10.28	93	20281	3.983 ng/ul#	93
27) 2,4-Dichlorophenol	10.52	162	13346	3.825 ng/ul	88
28) Naphthalene	10.93	128	46543	4.051 ng/ul	98
30) 4-Chloroaniline	11.03	127	16299	4.107 ng/ul#	89
31) Hexachlorobutadiene	11.22	225	10104	4.057 ng/ul#	86
32) Caprolactam	11.88	113	204	0.141 ng/ul	80
33) 4-Chloro-3-methylphenol	12.14	107	15307	3.734 ng/ul#	94
34) 2-Methylnaphthalene	12.53	142	35232	4.109 ng/ul	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.75	142	34153	4.084	ng/ul	94
37) 1,2,4,5-Tetrachlorobenzene	12.89	216	18843	3.814	ng/ul#	97
38) Hexachlorocyclopentadiene	12.88	237	9299	3.145	ng/ul#	89
39) 2,4,6-Trichlorophenol	13.13	196	10526	3.449	ng/ul	91
40) 2,4,5-Trichlorophenol	13.20	196	11758	3.717	ng/ul	98
41) 1,1'-Biphenyl	13.54	154	45329	3.993	ng/ul#	96
42) 2-Chloronaphthalene	13.58	162	36247	4.162	ng/ul	94
43) 2-Nitroaniline	13.77	65	9162	4.020	ng/ul#	72
45) Dimethylphthalate	14.15	163	45486	3.908	ng/ul	99
46) 2,6-Dinitrotoluene	14.27	165	6226	3.828	ng/ul	93
48) Acenaphthylene	14.42	152	53809	3.903	ng/ul	99
49) 3-Nitroaniline	14.59	138	6275	3.702	ng/ul#	92
50) Acenaphthene	14.76	153	38620	3.913	ng/ul	98
51) 2,4-Dinitrophenol	14.79	184	2357	2.849	ng/ul#	79
53) 4-Nitrophenol	14.88	109	6771	4.067	ng/ul	89
54) Dibenzofuran	15.09	168	55385	4.054	ng/ul	98
55) 2,4-Dinitrotoluene	15.05	165	8673	3.855	ng/ul	90
56) 2,3,4,6-Tetrachlorophenol	15.32	232	10762	3.633	ng/ul	98
57) Diethylphthalate	15.52	149	46448	3.932	ng/ul	96
59) Fluorene	15.75	166	43529	3.952	ng/ul	99
60) 4-Chlorophenyl-phenylether	15.74	204	24188	4.031	ng/ul	95
61) 4-Nitroaniline	15.75	138	6058	2.821	ng/ul#	86
64) 4,6-Dinitro-2-methylphenol	15.80	198	4986	3.726	ng/ul#	44
65) N-Nitrosodiphenylamine	15.95	169	40113	4.003	ng/ul	99
66) 4-Bromophenyl-phenylether	16.63	248	14975	3.785	ng/ul	95
67) Hexachlorobenzene	16.75	284	15738	3.986	ng/ul	96
68) Atrazine	16.89	200	15957	3.861	ng/ul	97
69) Pentachlorophenol	17.09	266	7628	3.186	ng/ul	92
70) Phenanthrene	17.49	178	75781	4.077	ng/ul	98
72) Anthracene	17.57	178	75505	4.037	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	13.50	216	20772	3.801	ng/uL	95
74) Pentachlorobenzene	15.01	250	18230	3.832	ng/uL	98
75) Carbazole	17.84	167	65616	4.173	ng/ul	97
76) Di-n-butylphthalate	18.41	149	78637	3.828	ng/ul	98
77) Fluoranthene	19.49	202	94240	4.373	ng/ul	97
80) Pyrene	19.85	202	91445	3.922	ng/ul#	94
81) Butylbenzylphthalate	20.75	149	32693	3.444	ng/ul#	85
82) 3,3'-Dichlorobenzidine	21.60	252	23972	3.024	ng/ul	95
83) Benzo(a)anthracene	21.69	228	95700	3.936	ng/ul	98
84) Bis(2-ethylhexyl)phthalate	21.62	149	49638	3.668	ng/ul	99
85) Chrysene	21.76	228	88340	3.961	ng/ul	98
87) Di-n-octyl phthalate	22.85	149	80388	3.821	ng/ul	100
88) Benzo(b)fluoranthene	23.91	252	85059	3.838	ng/ul#	94
89) Benzo(k)fluoranthene	23.98	252	87389	4.090	ng/ul#	97
91) Benzo(a)pyrene	24.79	252	82335	3.926	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	28.59	276	95280	3.815	ng/ul	99
93) Dibenzo(a,h)anthracene	28.67	278	45268	2.236	ng/ul	94
94) Benzo(g,h,i)perylene	29.72	276	39483	1.917	ng/ul	97

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

