

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082219\
 Data File : BM022240.D
 Acq On : 22 Aug 2019 16:07
 Operator : HP/JU
 Sample : SSTD00540
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD00540

Quant Time: Aug 22 18:25:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 22 18:08:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	22096	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	99391	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.22	164	70263	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.98	188	165263	20.00	ng/ul	0.00
77) Chrysene-d12	21.19	240	210355	20.00	ng/ul	0.00
85) Perylene-d12	23.39	264	250985	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.14	96	886	1.64	ng/uL	0.00
5) Phenol-d5	6.76	99	7590	4.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	5182	5.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.10	132	6895	4.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	6537	4.61	ng/ul	0.00
19) Nitrobenzene-d5	8.74	128	3048	4.10	ng/ul	0.00
22) 2-Nitrophenol-d4	9.45	143	3136	3.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	7253	4.27	ng/ul	0.01
29) 4-Chloroaniline-d4	10.52	131	7669	4.00	ng/ul	0.02
43) Dimethylphthalate-d6	13.65	166	29469	5.04	ng/ul	0.00
46) Acenaphthylene-d8	13.91	160	29680	4.46	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	1022	1.29	ng/ul	0.02
57) Fluorene-d10	15.22	176	24267	4.93	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.40	200	3027	2.63	ng/ul	0.02
70) Anthracene-d10	17.08	188	39948	4.67	ng/ul	0.00
78) Pyrene-d10	19.39	212	51543	4.58	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.24	264	61571	4.57	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.18	88	994	1.575	ng/uL 92
4) Benzaldehyde	6.75	77	4139	4.079	ng/ul 92
6) Phenol	6.79	94	8121	4.404	ng/ul 95
8) Bis(2-Chloroethyl)ether	7.02	93	6637	4.826	ng/ul 90
10) 2-Chlorophenol	7.13	128	6991	4.531	ng/ul 99
11) 2-Methylphenol	8.02	108	6317	4.570	ng/ul 86
14) Acetophenone	8.40	105	12390	5.146	ng/ul 96
15) N-Nitroso-di-n-propylamine	8.37	70	6376	5.454	ng/ul 98
16) 4-Methylphenol	8.35	108	7164	4.875	ng/ul 94
17) Hexachloroethane	8.61	117	3611	4.655	ng/ul 89
20) Nitrobenzene	8.78	77	9774	4.470	ng/ul 98
21) Isophorone	9.29	82	18318	5.042	ng/ul 99
23) 2-Nitrophenol	9.48	139	3523	3.754	ng/ul 95
24) 2,4-Dimethylphenol	9.54	107	9859	4.657	ng/ul 95
25) Bis(2-Chloroethoxy)methane	9.79	93	9531	4.662	ng/ul# 98
27) 2,4-Dichlorophenol	10.02	162	7315	4.374	ng/ul 94
28) Naphthalene	10.39	128	26425	4.871	ng/ul 99
30) 4-Chloroaniline	10.54	127	8342	4.189	ng/ul 96
31) Hexachlorobutadiene	10.64	225	8314	4.826	ng/ul 97
32) Caprolactam	11.37	113	750	1.657	ng/ul 88
33) 4-Chloro-3-methylphenol	11.67	107	8570	4.895	ng/ul 99
34) 2-Methylnaphthalene	12.01	142	19862	5.092	ng/ul 95
36) 1,2,4,5-Tetrachlorobenzene	12.39	216	13629	4.435	ng/ul 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) Hexachlorocyclopentadiene	12.33	237	2076	1.775	ng/ul#	93
38) 2,4,6-Trichlorophenol	12.65	196	6605	3.916	ng/ul	93
39) 2,4,5-Trichlorophenol	12.74	196	7150	3.862	ng/ul	95
40) 1,1'-Biphenyl	13.05	154	27082	4.644	ng/ul#	97
41) 2-Chloronaphthalene	13.09	162	21054	4.584	ng/ul	97
42) 2-Nitroaniline	13.33	65	5256	4.047	ng/ul	99
44) Dimethylphthalate	13.69	163	29912	5.016	ng/ul	99
45) 2,6-Dinitrotoluene	13.84	165	5460	4.892	ng/ul#	89
47) Acenaphthylene	13.94	152	32577	4.665	ng/ul	99
48) 3-Nitroaniline	14.19	138	3185	3.230	ng/ul#	95
49) Acenaphthene	14.28	153	23768	4.835	ng/ul	99
50) 2,4-Dinitrophenol	14.42	184	1254	2.074	ng/ul#	79
52) 4-Nitrophenol	14.52	109	2269	1.883	ng/ul	92
53) Dibenzofuran	14.63	168	34179	4.822	ng/ul	100
54) 2,4-Dinitrotoluene	14.64	165	7659	4.596	ng/ul#	95
55) 2,3,4,6-Tetrachlorophenol	14.87	232	7435	4.362	ng/ul#	97
56) Diethylphthalate	15.06	149	31423	5.079	ng/ul	97
58) Fluorene	15.28	166	27172	4.750	ng/ul	98
59) 4-Chlorophenyl-phenylether	15.27	204	16479	4.801	ng/ul	92
60) 4-Nitroaniline	15.36	138	1759	1.835	ng/ul#	76
63) 4,6-Dinitro-2-methylphenol	15.41	198	3675	2.963	ng/ul#	81
64) N-Nitrosodiphenylamine	15.50	169	23757	4.628	ng/ul	94
65) 4-Bromophenyl-phenylether	16.17	248	10647	4.555	ng/ul	93
66) Hexachlorobenzene	16.27	284	11876	4.601	ng/ul	96
67) Atrazine	16.47	200	10737	4.358	ng/ul	99
68) Pentachlorophenol	16.64	266	5061	3.457	ng/ul	90
69) Phenanthrene	17.02	178	45747	4.700	ng/ul	99
71) Anthracene	17.12	178	45705	4.551	ng/ul	99
72) 1,2,3,4-Tetrachlorobenzene	13.00	216	13775	4.151	ng/uL	98
73) Pentachlorobenzene	14.53	250	14786	4.445	ng/uL	92
74) Carbazole	17.41	167	38206	4.357	ng/ul	99
75) Di-n-butylphthalate	17.95	149	54094	5.088	ng/ul	97
76) Fluoranthene	19.05	202	61476	4.560	ng/ul#	96
79) Pyrene	19.42	202	63623	4.490	ng/ul	96
80) Butylbenzylphthalate	20.33	149	27076	4.854	ng/ul	98
81) 3,3'-Dichlorobenzidine	21.13	252	20275	3.899	ng/ul	95
82) Benzo(a)anthracene	21.17	228	73536	4.639	ng/ul	99
83) Bis(2-ethylhexyl)phthalate	21.10	149	42089	5.288	ng/ul	98
84) Chrysene	21.23	228	70284	4.725	ng/ul	99
86) Di-n-octyl phthalate	21.97	149	74735	5.053	ng/ul	100
87) Benzo(b)fluoranthene	22.73	252	75175	4.462	ng/ul	99
88) Benzo(k)fluoranthene	22.77	252	72886	4.408	ng/ul	98
90) Benzo(a)pyrene	23.29	252	72797	4.485	ng/ul	98
91) Indeno(1,2,3-cd)pyrene	25.58	276	85224	4.214	ng/ul	99
92) Dibenzo(a,h)anthracene	25.59	278	70068	4.139	ng/ul#	98
93) Benzo(g,h,i)perylene	26.26	276	71382	4.443	ng/ul	97

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

