

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
 Data File : BM020995.D
 Acq On : 18 Jun 2019 00:49
 Operator : HP/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02029

Quant Time: Jun 18 01:37:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 17 09:33:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	250144	20.00	ng/ul	0.00
18) Naphthalene-d8	10.58	136	1122793	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	749734	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1876344	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	1987488	20.00	ng/ul	0.00
85) Perylene-d12	23.63	264	1981045	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	38802	7.37	ng/uL	0.00
5) Phenol-d5	6.96	99	445239	20.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	268281	20.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	361625	20.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	380493	20.94	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	188715	20.70	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	214567	21.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	438945	21.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	496462	23.13	ng/ul	0.00
43) Dimethylphthalate-d6	13.84	166	1417521	20.92	ng/ul	-0.01
46) Acenaphthylene-d8	14.12	160	1729960	20.86	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	250857	22.77	ng/ul	0.00
57) Fluorene-d10	15.42	176	1238807	21.02	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.55	200	227385	19.71	ng/ul	0.00
70) Anthracene-d10	17.27	188	1981550	20.32	ng/ul	0.00
78) Pyrene-d10	19.57	212	2263764	18.98	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.49	264	2368798	20.27	ng/ul	-0.01

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.33	88	41633	7.810	ng/uL 96
4) Benzaldehyde	6.95	77	204538	20.606	ng/ul 98
6) Phenol	6.99	94	421210	20.453	ng/ul 100
8) Bis(2-Chloroethyl)ether	7.22	93	320668	20.402	ng/ul 98
10) 2-Chlorophenol	7.36	128	342289	20.531	ng/ul 99
11) 2-Methylphenol	8.23	108	329972	20.708	ng/ul 98
12) 2,2'-oxybis(1-Chloropropan	8.32	45	416198	20.737	ng/ul 98
14) Acetophenone	8.62	105	551405	21.080	ng/ul 98
15) N-Nitroso-di-n-propylamine	8.60	70	290207	21.317	ng/ul 98
16) 4-Methylphenol	8.56	108	363442	20.960	ng/ul 98
17) Hexachloroethane	8.86	117	138654	19.952	ng/ul 97
20) Nitrobenzene	8.99	77	416002	20.311	ng/ul 97
21) Isophorone	9.52	82	805939	20.894	ng/ul 99
23) 2-Nitrophenol	9.70	139	210146	20.651	ng/ul 96
24) 2,4-Dimethylphenol	9.76	107	428535	20.613	ng/ul 98
25) Bis(2-Chloroethoxy)methane	10.00	93	485285	20.416	ng/ul 99
27) 2,4-Dichlorophenol	10.23	162	392589	20.956	ng/ul 97
28) Naphthalene	10.63	128	1167002	20.244	ng/ul 100
30) 4-Chloroaniline	10.75	127	461403	22.856	ng/ul 97
31) Hexachlorobutadiene	10.90	225	250146	19.597	ng/ul 97
32) Caprolactam	11.53	113	122381	23.237	ng/ul 97
33) 4-Chloro-3-methylphenol	11.87	107	415590	21.759	ng/ul 97
34) 2-Methylnaphthalene	12.24	142	922510	20.871	ng/ul 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.61	216	526791	19.682	ng/ul	98
37) Hexachlorocyclopentadiene	12.58	237	276851	17.387	ng/ul	98
38) 2,4,6-Trichlorophenol	12.86	196	348074	20.927	ng/ul	97
39) 2,4,5-Trichlorophenol	12.93	196	377679	21.258	ng/ul	99
40) 1,1'-Biphenyl	13.26	154	1237811	20.014	ng/ul	99
41) 2-Chloronaphthalene	13.30	162	971923	19.891	ng/ul	99
42) 2-Nitroaniline	13.52	65	263624	22.200	ng/ul	99
44) Dimethylphthalate	13.89	163	1280542	20.765	ng/ul	99
45) 2,6-Dinitrotoluene	14.02	165	258467	21.960	ng/ul	98
47) Acenaphthylene	14.15	152	1472175	20.735	ng/ul	99
48) 3-Nitroaniline	14.35	138	239988	23.292	ng/ul	98
49) Acenaphthene	14.49	153	1069918	20.555	ng/ul	100
50) 2,4-Dinitrophenol	14.55	184	125134	21.163	ng/ul	99
52) 4-Nitrophenol	14.65	109	182465	22.387	ng/ul	99
53) Dibenzofuran	14.83	168	1534504	20.705	ng/ul	99
54) 2,4-Dinitrotoluene	14.80	165	392759	22.995	ng/ul	100
55) 2,3,4,6-Tetrachlorophenol	15.06	232	346797	22.384	ng/ul	98
56) Diethylphthalate	15.26	149	1287341	21.637	ng/ul	99
58) Fluorene	15.48	166	1269215	21.078	ng/ul	99
59) 4-Chlorophenyl-phenylether	15.47	204	675673	20.581	ng/ul	98
60) 4-Nitroaniline	15.51	138	261633	23.889	ng/ul	97
63) 4,6-Dinitro-2-methylphenol	15.56	198	230400	19.899	ng/ul	100
64) N-Nitrosodiphenylamine	15.69	169	1112919	19.585	ng/ul	100
65) 4-Bromophenyl-phenylether	16.37	248	445646	19.368	ng/ul	97
66) Hexachlorobenzene	16.47	284	519761	19.710	ng/ul	98
67) Atrazine	16.64	200	429609	20.838	ng/ul	97
68) Pentachlorophenol	16.82	266	287222	21.722	ng/ul	98
69) Phenanthrene	17.22	178	2091401	20.154	ng/ul	99
71) Anthracene	17.31	178	2145570	20.289	ng/ul	99
72) 1,2,3,4-Tetrachlorobenzene	13.22	216	529554	17.941	ng/ul	100
73) Pentachlorobenzene	14.75	250	567944	18.699	ng/ul	98
74) Carbazole	17.58	167	1879866	21.570	ng/ul	100
75) Di-n-butylphthalate	18.14	149	2248477	21.489	ng/ul	100
76) Fluoranthene	19.23	202	2627230	23.431	ng/ul	99
79) Pvrene	19.60	202	2654389	18.891	ng/ul	99
80) Butylbenzylphthalate	20.49	149	1103060	21.208	ng/ul	100
81) 3,3'-Dichlorobenzidine	21.27	252	931288	20.906	ng/ul	100
82) Benzo(a)anthracene	21.34	228	2724965	20.111	ng/ul	99
83) Bis(2-ethylhexyl)phthalate	21.26	149	1621427	21.895	ng/ul	99
84) Chrysene	21.39	228	2583673	20.360	ng/ul	100
86) Di-n-octyl phthalate	22.15	149	2777967	24.335	ng/ul	100
87) Benzo(b)fluoranthene	22.94	252	2691398	21.447	ng/ul	100
88) Benzo(k)fluoranthene	22.99	252	2470955	20.468	ng/ul	100
90) Benzo(a)pyrene	23.53	252	2365353	20.026	ng/ul	100
91) Indeno(1,2,3-cd)pyrene	25.93	276	2419348	17.395	ng/ul	99
92) Dibenzo(a,h)anthracene	25.94	278	2050531	17.525	ng/ul	99
93) Benzo(g,h,i)perylene	26.64	276	1910437	16.678	ng/ul	99

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

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