

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030421\
 Data File : BM028980.D
 Acq On : 04 Mar 2021 09:07
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02008

Quant Time: Mar 04 09:46:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 04 09:45:23 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	12467	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	53152	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.43	164	34093	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	72469	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	74172	20.00	ng/ul	0.00
86) Perylene-d12	23.53	264	69379	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	2264	8.18	ng/uL	0.00
5) Phenol-d5	6.95	99	18914	21.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	11108	21.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	16935	21.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	15998	22.07	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	7744	20.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	8761	20.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	16857	21.26	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	22410	24.89	ng/ul	0.00
44) Dimethylphthalate-d6	13.85	166	51707	22.06	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	63636	20.24	ng/ul	0.00
52) 4-Nitrophenol-d4	14.67	143	8316	20.39	ng/ul	0.00
58) Fluorene-d10	15.42	176	43128	21.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	8360	19.35	ng/ul	0.00
71) Anthracene-d10	17.27	188	67880	20.61	ng/ul	0.00
79) Pyrene-d10	19.56	212	73166	20.07	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.40	264	72553	20.59	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.16	88	2422	8.680	ng/uL 96
4) Benzaldehyde	6.91	77	11777	26.302	ng/ul 96
6) Phenol	6.97	94	19716	21.084	ng/ul 96
8) Bis(2-Chloroethyl)ether	7.20	93	15359	21.195	ng/ul 98
10) 2-Chlorophenol	7.33	128	17509	21.267	ng/ul 98
11) 2-Methylphenol	8.23	108	15282	21.653	ng/ul 100
12) 2,2'-oxybis(1-Chloropropan	8.30	45	26451	23.608	ng/ul 98
14) Acetophenone	8.60	105	26353	22.325	ng/ul 95
15) N-Nitroso-di-n-propylamine	8.59	70	12591	24.105	ng/ul 89
16) 4-Methylphenol	8.56	108	16629	22.148	ng/ul 97
17) Hexachloroethane	8.83	117	7682	21.826	ng/ul 95
20) Nitrobenzene	8.99	77	18939	21.099	ng/ul 98
21) Isophorone	9.50	82	33716	22.248	ng/ul 99
23) 2-Nitrophenol	9.69	139	9701	20.180	ng/ul 93
24) 2,4-Dimethylphenol	9.76	107	20466	22.130	ng/ul 93
25) Bis(2-Chloroethoxy)methane	9.99	93	21179	21.363	ng/ul# 98
27) 2,4-Dichlorophenol	10.22	162	16851	21.287	ng/ul 95
28) Naphthalene	10.62	128	55189	20.353	ng/ul 98
30) 4-Chloroaniline	10.74	127	22224	24.324	ng/ul 95
31) Hexachlorobutadiene	10.88	225	11067	21.439	ng/ul 94
32) Caprolactam	11.53	113	4952	23.384	ng/ul 88
33) 4-Chloro-3-methylphenol	11.88	107	18204	22.818	ng/ul 95
34) 2-Methylnaphthalene	12.23	142	40050	21.610	ng/ul 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.45	142	39289	22.022	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.60	216	20247	19.066	ng/ul	96
38) Hexachlorocyclopentadiene	12.57	237	6515	15.872	ng/ul	99
39) 2,4,6-Trichlorophenol	12.86	196	13550	20.192	ng/ul	91
40) 2,4,5-Trichlorophenol	12.93	196	14660	20.491	ng/ul	97
41) 1,1'-Biphenyl	13.26	154	51774	19.479	ng/ul	99
42) 2-Chloronaphthalene	13.30	162	41093	19.607	ng/ul	98
43) 2-Nitroaniline	13.52	65	11979	22.069	ng/ul	99
45) Dimethylphthalate	13.89	163	52696	21.555	ng/ul	100
46) 2,6-Dinitrotoluene	14.03	165	10745	22.318	ng/ul	100
48) Acenaphthylene	14.15	152	59997	20.107	ng/ul	98
49) 3-Nitroaniline	14.36	138	10883	23.233	ng/ul	99
50) Acenaphthene	14.49	153	43902	20.327	ng/ul	98
51) 2,4-Dinitrophenol	14.58	184	4872	18.035	ng/ul#	81
53) 4-Nitrophenol	14.68	109	8079	22.321	ng/ul	91
54) Dibenzofuran	14.83	168	62046	20.749	ng/ul	98
55) 2,4-Dinitrotoluene	14.82	165	15943	22.808	ng/ul#	88
56) 2,3,4,6-Tetrachlorophenol	15.06	232	11743	21.208	ng/ul#	90
57) Diethylphthalate	15.26	149	56390	22.816	ng/ul	99
59) Fluorene	15.48	166	51793	21.122	ng/ul	100
60) 4-Chlorophenyl-phenylether	15.47	204	25900	21.399	ng/ul	96
61) 4-Nitroaniline	15.52	138	11071	22.939	ng/ul	90
64) 4,6-Dinitro-2-methylphenol	15.59	198	8771	20.004	ng/ul	97
65) N-Nitrosodiphenylamine	15.69	169	44381	19.935	ng/ul	96
66) 4-Bromophenyl-phenylether	16.37	248	14953	20.273	ng/ul	94
67) Hexachlorobenzene	16.47	284	16701	20.162	ng/ul	100
68) Atrazine	16.64	200	16960	21.651	ng/ul	95
69) Pentachlorophenol	16.83	266	8676	19.505	ng/ul	94
70) Phenanthrene	17.22	178	80982	20.386	ng/ul	99
72) Anthracene	17.30	178	85033	20.771	ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	13.22	216	20454	17.989	ng/uL	96
74) Pentachlorobenzene	14.75	250	19956	19.055	ng/uL	95
75) Carbazole	17.58	167	75751	20.949	ng/ul	100
76) Di-n-butylphthalate	18.13	149	101221	23.286	ng/ul	99
77) Fluoranthene	19.23	202	95282	21.999	ng/ul	97
80) Pvrene	19.59	202	98914	20.357	ng/ul	99
81) Butylbenzylphthalate	20.48	149	46970	22.294	ng/ul	95
82) 3,3'-Dichlorobenzidine	21.26	252	33605	22.466	ng/ul	98
83) Benzo(a)anthracene	21.32	228	96681	20.857	ng/ul	100
84) Bis(2-ethylhexyl)phthalate	21.24	149	73610	23.388	ng/ul	98
85) Chrysene	21.37	228	94135	20.934	ng/ul	100
87) Di-n-octyl phthalate	22.11	149	127591	24.229	ng/ul	100
88) Benzo(b)fluoranthene	22.88	252	93628	20.768	ng/ul	98
89) Benzo(k)fluoranthene	22.92	252	93223	22.159	ng/ul	99
91) Benzo(a)pyrene	23.44	252	81687	20.861	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.74	276	92911	18.770	ng/ul	99
93) Dibenzo(a,h)anthracene	25.75	278	79627	18.847	ng/ul	97
94) Benzo(a,h,i)perylene	26.41	276	75956	18.370	ng/ul	97

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

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