

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011921\
 Data File : BM028457.D
 Acq On : 19 Jan 2021 13:21
 Operator : CG/JU
 Sample : SSTD16011
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD160011

Quant Time: Jan 19 13:53:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM011921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 19 13:19:35 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	30302	20.00	ng/ul	0.00
20) Naphthalene-d8	10.91	136	121932	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	68749	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.45	188	131393	20.00	ng/ul	0.00
79) Chrysene-d12	21.58	240	117999	20.00	ng/ul	0.01
88) Perylene-d12	23.89	264	116769	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	46303	61.97	ng/uL	0.00
4) Pyridine-d5	3.78	84	334180	156.97	ng/ul	0.00
7) Phenol-d5	7.25	99	405786	158.26	ng/ul	0.02
9) Bis-(2-Chloroethyl)ether-d	7.42	67	253956	160.40	ng/ul	0.02
11) 2-Chlorophenol-d4	7.62	132	353646	164.54	ng/ul	0.01
15) 4-Methylphenol-d8	8.81	113	331785	160.32	ng/ul	0.02
21) Nitrobenzene-d5	9.27	128	169557	171.68	ng/ul	0.01
24) 2-Nitrophenol-d4	9.99	143	190261	181.75	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.53	165	336200	170.67	ng/ul	0.02
31) 4-Chloroaniline-d4	11.05	131	444684	165.16	ng/ul	0.01
46) Dimethylphthalate-d6	14.14	166	878096	169.09	ng/ul	0.02
49) Acenaphthylene-d8	14.43	160	1101119	166.81	ng/ul	0.01
54) 4-Nitrophenol-d4	14.93	143	163876	163.77	ng/ul	0.03
60) Fluorene-d10	15.72	176	740049	160.92	ng/ul	0.01
65) 4,6-Dinitro-2-methylphenol	15.84	200	152789	179.58	ng/ul	0.02
73) Anthracene-d10	17.56	188	1062876	163.62	ng/ul	0.01
81) Pyrene-d10	19.82	212	1061444	156.62	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.76	264	1012275	161.47	ng/ul	0.02

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.36	88	47170	61.363	ng/uL 97
5) Pyridine	3.80	79	332618	155.260	ng/ul 96
6) Benzaldehyde	7.21	77	115620	116.353	ng/ul 99
8) Phenol	7.27	94	408211	157.110	ng/ul 99
10) Bis(2-Chloroethyl)ether	7.51	93	340056	160.507	ng/ul 97
12) 2-Chlorophenol	7.65	128	360753	162.839	ng/ul 99
13) 2-Methylphenol	8.53	108	315215	160.070	ng/ul 96
14) 2,2'-oxybis(1-Chloropropan	8.62	45	577826	165.088	ng/ul 99
16) Acetophenone	8.93	105	504024	159.626	ng/ul 98
17) N-Nitroso-di-n-propylamine	8.94	70	260751	167.500	ng/ul 95
18) 4-Methylphenol	8.89	108	336632	159.048	ng/ul 100
19) Hexachloroethane	9.17	117	163347	172.760	ng/ul 97
22) Nitrobenzene	9.32	77	395742	164.065	ng/ul 100
23) Isophorone	9.85	82	713267	175.002	ng/ul 98
25) 2-Nitrophenol	10.03	139	201551	174.312	ng/ul 96
26) 2,4-Dimethylphenol	10.08	107	375711	166.216	ng/ul 97
27) Bis(2-Chloroethoxy)methane	10.32	93	453101	168.023	ng/ul 100
29) 2,4-Dichlorophenol	10.56	162	325900	169.051	ng/ul 99
30) Naphthalene	10.96	128	1117273	162.007	ng/ul 99
32) 4-Chloroaniline	11.07	127	432130	164.411	ng/ul 98
33) Hexachlorobutadiene	11.23	225	216760	171.483	ng/ul 99
34) Caprolactam	11.90	113	57678	99.549	ng/ul 98
35) 4-Chloro-3-methylphenol	12.20	107	332548	166.835	ng/ul 98
36) 2-Methylnaphthalene	12.56	142	762256	165.338	ng/ul 100
37) 1-Methylnaphthalene	12.78	142	728545	164.092	ng/ul 100

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011921\
 Data File : BM028457.D
 Acq On : 19 Jan 2021 13:21
 Operator : CG/JU
 Sample : SSTD16011
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD160011

Quant Time: Jan 19 13:53:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM011921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 19 13:19:35 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
39) 1,2,4,5-Tetrachlorobenzene	12.93	216	374933	162.005	ng/ul	99
40) Hexachlorocyclopentadiene	12.90	237	230984	171.682	ng/ul	98
41) 2,4,6-Trichlorophenol	13.17	196	247304	171.886	ng/ul	99
42) 2,4,5-Trichlorophenol	13.24	196	261707	167.867	ng/ul	98
43) 1,1'-Biphenyl	13.57	154	962555	163.985	ng/ul	100
44) 2-Chloronaphthalene	13.62	162	746363	160.474	ng/ul	100
45) 2-Nitroaniline	13.82	65	223593	174.240	ng/ul	97
47) Dimethylphthalate	14.19	163	887275	164.604	ng/ul	98
48) 2,6-Dinitrotoluene	14.32	165	197629	184.722	ng/ul	96
50) Acenaphthylene	14.46	152	1114456	164.207	ng/ul	99
51) 3-Nitroaniline	14.64	138	164484	161.401	ng/ul	99
52) Acenaphthene	14.79	153	787196	161.547	ng/ul	98
53) 2,4-Dinitrophenol	14.85	184	118422	184.583	ng/ul	96
55) 4-Nitrophenol	14.95	109	125003	161.017	ng/ul	93
56) Dibenzofuran	15.13	168	1054852	160.657	ng/ul	99
57) 2,4-Dinitrotoluene	15.10	165	265539	176.803	ng/ul	96
58) 2,3,4,6-Tetrachlorophenol	15.35	232	220092	174.783	ng/ul	97
59) Diethylphthalate	15.54	149	928841	173.805	ng/ul	99
61) Fluorene	15.77	166	878727	160.724	ng/ul	99
62) 4-Chlorophenyl-phenylether	15.76	204	432778	163.326	ng/ul	99
63) 4-Nitroaniline	15.80	138	140057	157.371	ng/ul	98
66) 4,6-Dinitro-2-methylphenol	15.86	198	160348	174.619	ng/ul#	95
67) N-Nitrosodiphenylamine	15.97	169	728311	168.006	ng/ul	98
68) 4-Bromophenyl-phenylether	16.64	248	264871	173.589	ng/ul	100
69) Hexachlorobenzene	16.77	284	286584	163.455	ng/ul	98
70) Atrazine	16.92	200	264254	175.705	ng/ul	100
71) Pentachlorophenol	17.10	266	179034	176.556	ng/ul	99
72) Phenanthrene	17.50	178	1286935	160.780	ng/ul	99
74) Anthracene	17.59	178	1325941	162.952	ng/ul	99
75) 1,2,3,4-Tetrachlorobenzene	13.53	216	366446	166.399	ng/uL	99
76) Pentachlorobenzene	15.05	250	351250	164.173	ng/uL	98
77) Carbazole	17.86	167	1121850	160.460	ng/ul	100
78) Di-n-butylphthalate	18.40	149	1601833	197.550	ng/ul	100
80) Fluoranthene	19.49	202	1395188	156.765	ng/ul	97
82) Pyrene	19.85	202	1413556	153.535	ng/ul	98
83) Butylbenzylphthalate	20.72	149	685959	201.159	ng/ul	98
84) 3,3'-Dichlorobenzidine	21.49	252	332281	134.244	ng/ul	99
85) Benzo(a)anthracene	21.56	228	1251479	158.474	ng/ul	99
86) Bis(2-ethylhexyl)phthalate	21.47	149	1053968	225.152	ng/ul	99
87) Chrysene	21.62	228	1204237	154.041	ng/ul	99
89) Di-n-octyl phthalate	22.37	149	1727353	206.370	ng/ul	100
90) Benzo(b)fluoranthene	23.20	252	1239588	157.966	ng/ul	99
91) Benzo(k)fluoranthene	23.25	252	1194200	154.000	ng/ul	99
93) Benzo(a)pyrene	23.81	252	1126478	158.428	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.29	276	1353233	162.343	ng/ul	95
95) Dibenzo(a,h)anthracene	26.30	278	1134632	160.495	ng/ul	98
96) Benzo(g,h,i)perylene	27.02	276	1151836	162.687	ng/ul	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011921\
 Data File : BM028457.D
 Acq On : 19 Jan 2021 13:21
 Operator : CG/JU
 Sample : SSTD16011
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD160011

Quant Time: Jan 19 13:53:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM011921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 19 13:19:35 2021
 Response via : Initial Calibration

