

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032921\
 Data File : BP005070.D
 Acq On : 29 Mar 2021 10:28
 Operator : CG/JU
 Sample : SSTD00576
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTD00576

Quant Time: Mar 29 12:12:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP032921MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 29 12:06:34 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	24524	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	96014	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	62165	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.13	188	137289	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	142786	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	148460	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	1454	2.18	ng/uL	0.00
5) Phenol-d5	6.89	99	10007	4.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	5731	5.04	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	6859	4.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	7418	4.53	ng/ul	0.00
19) Nitrobenzene-d5	8.88	128	3369	4.73	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	3563	4.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	6953	4.49	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	7481	4.41	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	21419	4.76	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	26368	4.74	ng/ul	0.00
52) 4-Nitrophenol-d4	14.59	143	1947	2.53	ng/ul	0.00
58) Fluorene-d10	15.37	176	19174	4.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	3106	3.37	ng/ul	0.00
71) Anthracene-d10	17.13	188	137289	22.11	ng/ul	-0.10
79) Pyrene-d10	19.47	212	33583	4.90	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	36039	4.66	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.26	88	1359	1.950	ng/uL 90
4) Benzaldehyde	6.86	77	6316	5.889	ng/ul 86
6) Phenol	6.92	94	9994	4.382	ng/ul 92
8) Bis(2-Chloroethyl)ether	7.15	93	8186	4.876	ng/ul 96
10) 2-Chlorophenol	7.28	128	7724	4.899	ng/ul 96
11) 2-Methylphenol	8.16	108	7326	4.402	ng/ul 94
12) 2,2'-oxybis(1-Chloropropan	8.26	45	6208	4.961	ng/ul 93
14) Acetophenone	8.54	105	12754	4.699	ng/ul 93
15) N-Nitroso-di-n-propylamine	8.52	70	6539	4.786	ng/ul 93
16) 4-Methylphenol	8.49	108	7883	4.426	ng/ul 96
17) Hexachloroethane	8.79	117	3600	5.074	ng/ul 93
20) Nitrobenzene	8.92	77	9795	4.696	ng/ul# 83
21) Isophorone	9.44	82	16210	4.488	ng/ul# 96
23) 2-Nitrophenol	9.62	139	3908	4.539	ng/ul 97
24) 2,4-Dimethylphenol	9.69	107	9546	4.678	ng/ul 93
25) Bis(2-Chloroethoxy)methane	9.93	93	10250	4.731	ng/ul 98
27) 2,4-Dichlorophenol	10.16	162	7310	4.769	ng/ul 95
28) Naphthalene	10.55	128	25219	4.992	ng/ul 98
30) 4-Chloroaniline	10.68	127	7484	4.226	ng/ul 90
31) Hexachlorobutadiene	10.84	225	5483	5.055	ng/ul 93
32) Caprolactam	11.46	113	1661	3.185	ng/ul# 77
33) 4-Chloro-3-methylphenol	11.80	107	7842	4.312	ng/ul# 79
34) 2-Methylnaphthalene	12.18	142	18272	5.080	ng/ul 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.39	142	16610	4.806	ng/ul	97
37) 1,2,4,5-Tetrachlorobenzene	12.55	216	10344	5.058	ng/ul#	93
38) Hexachlorocyclopentadiene	12.52	237	4883	3.932	ng/ul#	88
39) 2,4,6-Trichlorophenol	12.79	196	5879	4.595	ng/ul	98
40) 2,4,5-Trichlorophenol	12.87	196	6073	4.525	ng/ul	98
41) 1,1'-Biphenyl	13.20	154	23194	4.995	ng/ul	92
42) 2-Chloronaphthalene	13.24	162	17977	4.847	ng/ul#	91
43) 2-Nitroaniline	13.45	65	4611	4.149	ng/ul#	90
45) Dimethylphthalate	13.83	163	22632	4.938	ng/ul	95
46) 2,6-Dinitrotoluene	13.95	165	4200	4.376	ng/ul	93
48) Acenaphthylene	14.09	152	25591	4.794	ng/ul	99
49) 3-Nitroaniline	14.28	138	2652	3.082	ng/ul	88
50) Acenaphthene	14.43	153	19127	5.050	ng/ul	95
51) 2,4-Dinitrophenol	14.49	184	852	1.419	ng/ul#	90
53) 4-Nitrophenol	14.60	109	1885	2.392	ng/ul	90
54) Dibenzofuran	14.77	168	27616	4.985	ng/ul	95
55) 2,4-Dinitrotoluene	14.74	165	5878	4.316	ng/ul#	99
56) 2,3,4,6-Tetrachlorophenol	15.00	232	5799	4.624	ng/ul#	98
57) Diethylphthalate	15.20	149	22080	4.882	ng/ul	99
59) Fluorene	15.42	166	22946	4.956	ng/ul	99
60) 4-Chlorophenyl-phenylether	15.42	204	12091	4.882	ng/ul	96
61) 4-Nitroaniline	15.45	138	2112	2.275	ng/ul#	86
64) 4,6-Dinitro-2-methylphenol	15.50	198	3143	3.402	ng/ul#	86
65) N-Nitrosodiphenylamine	15.64	169	18591	4.625	ng/ul	98
66) 4-Bromophenyl-phenylether	16.32	248	7339	5.003	ng/ul	95
67) Hexachlorobenzene	16.43	284	8629	5.078	ng/ul	98
68) Atrazine	16.60	200	6833	4.369	ng/ul	96
69) Pentachlorophenol	16.78	266	4294	3.978	ng/ul	97
70) Phenanthrene	17.17	178	36819	4.924	ng/ul	96
72) Anthracene	17.26	178	36383	4.780	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	13.16	216	10138	4.749	ng/uL	95
74) Pentachlorobenzene	14.69	250	10000	4.898	ng/uL	99
75) Carbazole	17.54	167	29716	4.552	ng/ul	96
76) Di-n-butylphthalate	18.10	149	32310	4.295	ng/ul	98
77) Fluoranthene	19.15	202	41492	4.787	ng/ul	98
80) Pvrene	19.50	202	44151	5.014	ng/ul	99
81) Butylbenzylphthalate	20.38	149	14666	4.536	ng/ul	99
82) 3,3'-Dichlorobenzidine	21.15	252	9402	3.176	ng/ul	97
83) Benzo(a)anthracene	21.21	228	44413	5.034	ng/ul	95
84) Bis(2-ethylhexyl)phthalate	21.14	149	20860	4.289	ng/ul	96
85) Chrysene	21.26	228	43600	5.099	ng/ul	96
87) Di-n-octyl phthalate	22.03	149	36627	3.860	ng/ul	100
88) Benzo(b)fluoranthene	22.82	252	47023	4.798	ng/ul	100
89) Benzo(k)fluoranthene	22.86	252	44015	4.652	ng/ul	95
91) Benzo(a)pyrene	23.41	252	39614	4.705	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	25.84	276	49445	4.592	ng/ul	98
93) Dibenzo(a,h)anthracene	25.86	278	40671	4.472	ng/ul	97
94) Benzo(a,h,i)perylene	26.57	276	43511	4.855	ng/ul	99

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

