

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020821\
 Data File : BP004690.D
 Acq On : 08 Feb 2021 13:11
 Operator : CG/JU
 Sample : SSTD04081
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTD04081

Quant Time: Feb 08 13:47:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 08 13:10:34 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	49478	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	202929	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	135604	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	301287	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	326506	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	318141	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	19105	19.95	ng/uL	0.00
5) Phenol-d5	6.94	99	160315	45.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	83234	40.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	122104	48.06	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	127383	46.38	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	57809	42.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	65964	38.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	128265	32.77	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	154194	46.96	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	393726	39.17	ng/ul	0.00
47) Acenaphthylene-d8	14.11	160	482733	40.20	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	70607	47.47	ng/ul	0.00
58) Fluorene-d10	15.42	176	333698	36.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	72732	39.95	ng/ul	0.00
71) Anthracene-d10	17.27	188	533843	39.83	ng/ul	0.00
79) Pyrene-d10	19.51	212	621842	38.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	656423	39.87	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.29	88	20225	20.468	ng/uL 91
4) Benzaldehyde	6.92	77	100675	46.573	ng/ul 98
6) Phenol	6.96	94	167578	44.513	ng/ul 99
8) Bis(2-Chloroethyl)ether	7.20	93	122609	45.200	ng/ul 98
10) 2-Chlorophenol	7.34	128	128835	47.848	ng/ul 97
11) 2-Methylphenol	8.21	108	125655	44.497	ng/ul 89
12) 2,2'-oxybis(1-Chloropropan	8.30	45	95502	33.618	ng/ul 98
14) Acetophenone	8.59	105	209671	43.889	ng/ul 99
15) N-Nitroso-di-n-propylamine	8.59	70	104156	41.415	ng/ul 98
16) 4-Methylphenol	8.55	108	134320	45.117	ng/ul 98
17) Hexachloroethane	8.85	117	55643	44.790	ng/ul 97
20) Nitrobenzene	8.98	77	161043	36.594	ng/ul 98
21) Isophorone	9.50	82	281491	37.183	ng/ul 99
23) 2-Nitrophenol	9.68	139	71878	41.206	ng/ul 97
24) 2,4-Dimethylphenol	9.74	107	167599	39.428	ng/ul 99
25) Bis(2-Chloroethoxy)methane	9.98	93	161454	39.088	ng/ul 98
27) 2,4-Dichlorophenol	10.21	162	126659	32.915	ng/ul 97
28) Naphthalene	10.62	128	411265	41.460	ng/ul 97
30) 4-Chloroaniline	10.73	127	161798	47.480	ng/ul 98
31) Hexachlorobutadiene	10.90	225	90312	26.544	ng/ul 97
32) Caprolactam	11.49	113	25814	26.482	ng/ul 98
33) 4-Chloro-3-methylphenol	11.85	107	152762	40.438	ng/ul 99
34) 2-Methylnaphthalene	12.23	142	301206	37.086	ng/ul 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.45	142	288135	36.426	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.60	216	172013	33.876	ng/ul	98
38) Hexachlorocyclopentadiene	12.59	237	109825	30.207	ng/ul	97
39) 2,4,6-Trichlorophenol	12.85	196	107056	35.843	ng/ul	95
40) 2,4,5-Trichlorophenol	12.91	196	118993	36.950	ng/ul	97
41) 1,1'-Biphenyl	13.25	154	387814	40.577	ng/ul	99
42) 2-Chloronaphthalene	13.29	162	307368	39.463	ng/ul	97
43) 2-Nitroaniline	13.50	65	95580	45.710	ng/ul	96
45) Dimethylphthalate	13.88	163	405018	39.707	ng/ul	100
46) 2,6-Dinitrotoluene	13.99	165	82917	41.564	ng/ul	96
48) Acenaphthylene	14.14	152	457460	42.446	ng/ul	100
49) 3-Nitroaniline	14.32	138	76229	52.226	ng/ul#	95
50) Acenaphthene	14.49	153	329645	41.798	ng/ul	98
51) 2,4-Dinitrophenol	14.53	184	49139	35.945	ng/ul	97
53) 4-Nitrophenol	14.63	109	80367	41.319	ng/ul	94
54) Dibenzofuran	14.82	168	467336	38.104	ng/ul	99
55) 2,4-Dinitrotoluene	14.78	165	119716	42.541	ng/ul	96
56) 2,3,4,6-Tetrachlorophenol	15.05	232	105370	33.603	ng/ul	96
57) Diethylphthalate	15.25	149	414138	42.746	ng/ul	100
59) Fluorene	15.47	166	397911	38.614	ng/ul	99
60) 4-Chlorophenyl-phenylether	15.47	204	212901	34.386	ng/ul	96
61) 4-Nitroaniline	15.49	138	85103	50.415	ng/ul	94
64) 4,6-Dinitro-2-methylphenol	15.55	198	73704	39.689	ng/ul#	91
65) N-Nitrosodiphenylamine	15.68	169	337976	42.603	ng/ul	97
66) 4-Bromophenyl-phenylether	16.36	248	129276	35.076	ng/ul	98
67) Hexachlorobenzene	16.47	284	141307	34.113	ng/ul	100
68) Atrazine	16.64	200	137930	37.832	ng/ul	99
69) Pentachlorophenol	16.82	266	85749	33.195	ng/ul	95
70) Phenanthrene	17.22	178	634804	40.796	ng/ul	100
72) Anthracene	17.30	178	648805	40.857	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	13.21	216	172954	37.313	ng/uL	98
74) Pentachlorobenzene	14.74	250	167928	34.590	ng/uL	96
75) Carbazole	17.57	167	565325	42.469	ng/ul	100
76) Di-n-butylphthalate	18.14	149	688165	47.453	ng/ul	99
77) Fluoranthene	19.19	202	769890	38.053	ng/ul	98
80) Pvrene	19.54	202	795949	39.755	ng/ul	99
81) Butylbenzylphthalate	20.42	149	307725	50.603	ng/ul	93
82) 3,3'-Dichlorobenzidine	21.17	252	287210	41.346	ng/ul#	98
83) Benzo(a)anthracene	21.24	228	791628	40.197	ng/ul	98
84) Bis(2-ethylhexyl)phthalate	21.17	149	465878	50.434	ng/ul	99
85) Chrysene	21.29	228	759971	39.740	ng/ul	99
87) Di-n-octyl phthalate	22.06	149	772601	48.870	ng/ul	100
88) Benzo(b)fluoranthene	22.86	252	823757	40.647	ng/ul	99
89) Benzo(k)fluoranthene	22.91	252	770392	39.197	ng/ul	100
91) Benzo(a)pyrene	23.46	252	708124	40.826	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.93	276	898375	40.294	ng/ul	97
93) Dibenzo(a,h)anthracene	25.94	278	758491	40.062	ng/ul	98
94) Benzo(a,h,i)perylene	26.66	276	731028	39.590	ng/ul	99

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

