

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121420\
 Data File : BP004297.D
 Acq On : 15 Dec 2020 10:42
 Operator : CG/JU
 Sample : L5001-23
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A54K6

Quant Time: Dec 15 11:12:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 15 09:00:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	278842	20.00	ng/ul	0.01
18) Naphthalene-d8	10.65	136	1192513	20.00	ng/ul	0.01
36) Acenaphthene-d10	14.49	164	737247	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.26	188	1583771	20.00	ng/ul	0.01
78) Chrysene-d12	21.34	240	1547769	20.00	ng/ul	0.00
86) Perylene-d12	23.72	264	1666376	20.00	ng/ul	0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	30413	3.83	ng/uL	0.00
5) Phenol-d5	7.01	99	775630	33.62	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.18	67	416182	28.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.38	132	604115	34.51	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	606704	34.81	ng/ul	0.01
19) Nitrobenzene-d5	9.00	128	309929	31.73	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	337565	41.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.26	165	655738	37.02	ng/ul	0.01
29) 4-Chloroaniline-d4	10.77	131	813677	37.06	ng/ul	0.01
44) Dimethylphthalate-d6	13.90	166	1941200	34.45	ng/ul	0.00
47) Acenaphthylene-d8	14.19	160	2411098	33.53	ng/ul	0.01
52) 4-Nitrophenol-d4	14.69	143	344087	36.15	ng/ul	0.00
58) Fluorene-d10	15.49	176	1665888	32.64	ng/ul	0.01
63) 4,6-Dinitro-2-methylphenol	15.62	200	188966	24.87	ng/ul	0.00
71) Anthracene-d10	17.36	188	2747948	35.60	ng/ul	0.01
79) Pyrene-d10	19.60	212	3133505	38.86	ng/ul	0.01
90) Benzo(a)pyrene-d12	23.57	264	3296524	37.55	ng/ul	0.02

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.36	88	129292	15.674	ng/uL 97
8) Bis(2-Chloroethyl)ether	7.27	93	803971	40.494	ng/ul 97
10) 2-Chlorophenol	7.41	128	752038	40.975	ng/ul 97
12) 2,2'-oxybis(1-Chloropropan	8.38	45	712461	30.020	ng/ul 98
15) N-Nitroso-di-n-propylamine	8.65	70	313455	23.479	ng/ul 97
17) Hexachloroethane	8.92	117	196778	23.345	ng/ul 95
20) Nitrobenzene	9.05	77	762949	31.637	ng/ul 99
25) Bis(2-Chloroethoxy)methane	10.05	93	765361	27.603	ng/ul 99
27) 2,4-Dichlorophenol	10.29	162	673777	37.489	ng/ul 99
28) Naphthalene	10.69	128	1624105	24.023	ng/ul 99
31) Hexachlorobutadiene	10.98	225	266945	19.227	ng/ul 99
37) 1,2,4,5-Tetrachlorobenzene	12.68	216	349638	13.021	ng/ul 99
38) Hexachlorocyclopentadiene	12.66	237	256151	16.626	ng/ul 99
39) 2,4,6-Trichlorophenol	12.92	196	775314	59.827	ng/ul 99
40) 2,4,5-Trichlorophenol	12.99	196	517141	35.690	ng/ul 97
42) 2-Chloronaphthalene	13.36	162	1303876	26.450	ng/ul 98
45) Dimethylphthalate	13.95	163	2074761	36.150	ng/ul 100
46) 2,6-Dinitrotoluene	14.07	165	239558	22.415	ng/ul 91
48) Acenaphthylene	14.22	152	1709890	23.958	ng/ul 100
50) Acenaphthene	14.56	153	1013511	20.056	ng/ul 98
54) Dibenzofuran	14.89	168	1631068	22.908	ng/ul 99
55) 2,4-Dinitrotoluene	14.86	165	276944	18.259	ng/ul 98
57) Diethylphthalate	15.32	149	1509318	28.019	ng/ul 99

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59) Fluorene	15.55	166	1373939	24.126	ng/ul	99
60) 4-Chlorophenyl-phenylether	15.54	204	425589	14.096	ng/ul	97
66) 4-Bromophenyl-phenylether	16.44	248	434662	23.599	ng/ul	99
67) Hexachlorobenzene	16.56	284	356246	16.133	ng/ul	97
69) Pentachlorophenol	16.90	266	424901	41.832	ng/ul	98
72) Anthracene	17.39	178	3540135	38.296	ng/ul	100
75) Carbazole	17.66	167	2700619	35.900	ng/ul	99
76) Di-n-butylphthalate	18.22	149	2970897	37.067	ng/ul	99
81) Butylbenzylphthalate	20.50	149	1521593	52.803	ng/ul	96
83) Benzo(a)anthracene	21.33	228	3118277	30.565	ng/ul	99
84) Bis(2-ethylhexyl)phthalate	21.26	149	1353676	29.008	ng/ul	100
87) Di-n-octyl phthalate	22.17	149	4299487	54.539	ng/ul	100
88) Benzo(b)fluoranthene	23.00	252	1632206	15.445	ng/ul	100
89) Benzo(k)fluoranthene	23.04	252	4345844	40.043	ng/ul	100
94) Benzo(a,h,i)perylene	26.90	276	2693067	25.957	ng/ul	96

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

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