

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
 Data File : BN010653.D
 Acq On : 29 Apr 2020 13:10
 Operator : CG/JU
 Sample : L2321-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2M15

Quant Time: Apr 29 13:51:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 29 11:04:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	452375	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	1989350	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.21	164	1262838	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.96	188	2783633	20.00	ng/ul	0.00
78) Chrysene-d12	21.17	240	2487540	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	2271896	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	27469	3.18	ng/uL	0.00
5) Phenol-d5	6.78	99	866815	24.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	436081	24.14	ng/ul	0.00
9) 2-Chlorophenol-d4	7.12	132	707144	23.22	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	660759	21.94	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	368043	24.01	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	402412	23.56	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	770919	23.02	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	955315	24.88	ng/ul	0.00
44) Dimethylphthalate-d6	13.63	166	2535873	27.06	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	2921823	25.38	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	432482	23.71	ng/ul	0.00
58) Fluorene-d10	15.20	176	2139514	25.44	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.33	200	286918	19.73	ng/ul	0.00
71) Anthracene-d10	17.06	188	3787437	29.38	ng/ul	0.00
79) Pyrene-d10	19.36	212	4478424	31.89	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.20	264	3722781	31.59	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.80	94	1763527	47.698	ng/ul 99
20) Nitrobenzene	8.76	77	1427386	39.866	ng/ul 98
21) Isophorone	9.29	82	3228427	48.885	ng/ul 99
28) Naphthalene	10.39	128	5594314	52.326	ng/ul 99
34) 2-Methylnaphthalene	12.00	142	3217306	42.579	ng/ul 100
39) 2,4,6-Trichlorophenol	12.63	196	1453100	53.260	ng/ul 98
41) 1,1'-Biphenyl	13.03	154	3487795	35.961	ng/ul 100
48) Acenaphthylene	13.92	152	2764509	22.757	ng/ul 99
54) Dibenzofuran	14.61	168	5400001	46.402	ng/ul 99
57) Diethylphthalate	15.05	149	3834670	39.207	ng/ul 98
67) Hexachlorobenzene	16.27	284	1151992	33.324	ng/ul 96
69) Pentachlorophenol	16.62	266	1278709	57.939	ng/ul 96
72) Anthracene	17.09	178	6876283	43.882	ng/ul 98
75) Carbazole	17.36	167	7953974	61.796	ng/ul 98
77) Fluoranthene	19.03	202	8297614	49.403	ng/ul 96
83) Benzo(a)anthracene	21.15	228	6173950	35.778	ng/ul 98
84) Bis(2-ethylhexyl)phthalate	21.11	149	5392132	48.157	ng/ul 100
87) Di-n-octyl phthalate	21.98	149	9603133	59.426	ng/ul 100
92) Indeno(1,2,3-cd)pyrene	25.49	276	7172442	48.379	ng/ul 96
93) Dibenz(a,h)anthracene	25.49	278	6270680	50.361	ng/ul 98
94) Benzo(a,h,i)perylene	26.13	276	6307858	51.412	ng/ul 98

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

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