

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025841.D
 Acq On : 27 Mar 2020 17:04
 Operator : CG/JU
 Sample : L1968-01DL 20X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0DC7DL

Quant Time: Mar 29 00:22:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 27 23:44:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	179778	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	731979	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.20	164	449998	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.95	188	974570	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	1051161	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	1150948	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.75	99	2979	0.21	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.91	67	9695	1.20	ng/ul	0.00
9) 2-Chlorophenol-d4	7.10	132	9954	0.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	5813	0.50	ng/ul	0.00
19) Nitrobenzene-d5	8.71	128	6388	1.17	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	5779	1.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	10878	0.92	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.62	166	43402	1.27	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	48714	1.19	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.20	176	40167	1.34	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	17.04	188	64045	1.41	ng/ul	0.00
79) Pyrene-d10	19.34	212	64913	1.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	80101	1.36	ng/ul	0.00

Target Compounds

					Ovalue
20) Nitrobenzene	8.75	77	298866	22.549 ng/ul	99
37) 1,2,4,5-Tetrachlorobenzene	12.38	216	18314	1.287 ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	12.99	216	25151	1.684 ng/uL	96

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

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