

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM010719\
 Data File : BM018446.D
 Acq On : 07 Jan 2019 16:14
 Operator : JU/SJ
 Sample : K1048-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0950

Quant Time: Jan 08 09:18:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM010719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 07 23:55:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	386266	20.00	ng/ul	0.04
18) Naphthalene-d8	10.57	136	1609363	20.00	ng/ul	0.02
35) Acenaphthene-d10	14.41	164	867166	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1959801	20.00	ng/ul	0.00
77) Chrysene-d12	21.35	240	2145571	20.00	ng/ul	0.00
85) Perylene-d12	23.64	264	2217165	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	34544	4.17	ng/uL	0.00
5) Phenol-d5	6.95	99	207748	6.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	527658	30.78	ng/ul	0.01
9) 2-Chlorophenol-d4	7.31	132	694843	26.05	ng/ul	0.01
13) 4-Methylphenol-d8	8.48	113	409087	16.56	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	412224	33.46	ng/ul	0.01
22) 2-Nitrophenol-d4	9.65	143	448291	35.65	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	765753	30.17	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	21861	1.11	ng/ul	0.02
43) Dimethylphthalate-d6	13.82	166	2561535	34.42	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	3230107	34.98	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	84166	6.55	ng/ul	0.00
57) Fluorene-d10	15.40	176	2217012	34.72	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.52	200	334374	31.16	ng/ul	0.00
70) Anthracene-d10	17.26	188	3423564	35.91	ng/ul	0.00
78) Pyrene-d10	19.56	212	3812139	34.95	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.49	264	4310242	35.75	ng/ul	0.00

Target Compounds

					Ovalue	
10) 2-Chlorophenol	7.34	128	46274	1.750	ng/ul	95
20) Nitrobenzene	8.99	77	3915706	145.656	ng/ul	98
23) 2-Nitrophenol	9.68	139	35496	2.611	ng/ul	98
36) 1,2,4,5-Tetrachlorobenzene	12.59	216	1234719	41.914	ng/ul	99
72) 1,2,3,4-Tetrachlorobenzene	13.20	216	3497749	123.126	ng/uL	99
73) Pentachlorobenzene	14.72	250	84283	3.099	ng/uL	95

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

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