

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102418\
 Data File : BM017319.D
 Acq On : 24 Oct 2018 16:12
 Operator : MJ/SJ
 Sample : J5567-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ1

Quant Time: Oct 25 01:51:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 25 01:24:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	251972	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1133102	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	673213	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	1612295	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	1978800	20.00	ng/ul	0.00
86) Perylene-d12	23.46	264	2230731	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	17911	3.27	ng/uL	0.00
5) Phenol-d5	6.84	99	399004	17.33	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	258942	19.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	331505	18.58	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	339096	18.63	ng/ul	0.00
19) Nitrobenzene-d5	8.81	128	171635	19.38	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	185922	18.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	340246	18.44	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	329818	21.90	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	1212655	21.00	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	1433709	20.45	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	180417	16.73	ng/ul	0.00
58) Fluorene-d10	15.30	176	1049418	20.93	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.42	200	169816	15.12	ng/ul	0.00
71) Anthracene-d10	17.15	188	1694555	21.23	ng/ul	0.00
79) Pyrene-d10	19.45	212	1987597	21.68	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.32	264	2575356	21.86	ng/ul	0.00

Target Compounds

					Qvalue
6) Phenol	6.86	94	119021	5.208 ng/ul	99
45) Dimethylphthalate	13.76	163	156162	2.792 ng/ul#	99

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

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