

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061818\
 Data File : BM015713.D
 Acq On : 19 Jun 2018 04:40
 Operator : SJ/JU
 Sample : J3527-13 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jun 19 05:12:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060818MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 19 05:11:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	91711	20.00	ng/ul	0.00
18) Naphthalene-d8	11.15	136	396864	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.94	164	223236	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.67	188	526401	20.00	ng/ul	0.00
77) Chrysene-d12	21.77	240	636459	20.00	ng/ul	0.00
85) Perylene-d12	24.18	264	696192	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.55	96	523	0.28	ng/uL	0.01
5) Phenol-d5	7.50	99	110	0.01	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.66	67	2644	0.53	ng/ul	0.02
9) 2-Chlorophenol-d4	7.87	132	3378	0.53	ng/ul	0.02
13) 4-Methylphenol-d8	9.07	113	1389	0.19	ng/ul	0.03
19) Nitrobenzene-d5	9.53	128	855	0.32	ng/ul	0.02
22) 2-Nitrophenol-d4	10.25	143	629	0.22	ng/ul	0.02
26) 2,4-Dichlorophenol-d3	10.80	165	1898	0.32	ng/ul	0.03
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	14.34	166	11357	0.61	ng/ul	0.00
46) Acenaphthylene-d8	14.65	160	11222	0.56	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.93	176	10913	0.70	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.77	188	15604	0.65	ng/ul	0.00
78) Pyrene-d10	20.02	212	18052	0.66	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.03	264	22532	0.65	ng/ul	0.00

Target Compounds Ovalue

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

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