

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
 Data File : BM016969.D
 Acq On : 03 Oct 2018 08:43
 Operator : MJ/SJ
 Sample : J5126-13
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BA1

Quant Time: Oct 03 11:36:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 02 20:15:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	344370	20.00	ng/ul	0.00
18) Naphthalene-d8	10.49	136	1503642	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	816859	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.10	188	1525165	20.00	ng/ul	0.00
78) Chrysene-d12	21.29	240	1216625	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	1215051	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	0.00	99	0d	0.00	ng/ul
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	ng/ul
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul
44) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul
58) Fluorene-d10	0.00	176	0d	0.00	ng/ul
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul
79) Pyrene-d10	0.00	212	0d	0.00	ng/ul
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds

				Ovalue	
73) 1,2,3,4-Tetrachlorobenzene	13.14	216	46087	1.826	ng/uL 95
77) Fluoranthene	19.16	202	124400	1.281	ng/ul# 89
85) Chrysene	21.33	228	76695	1.092	ng/ul 96
88) Benzo(b)fluoranthene	22.85	252	85040	1.169	ng/ul# 65

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

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