

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101318\
 Data File : BM017158.D
 Acq On : 13 Oct 2018 21:46
 Operator : MJ/SJ
 Sample : J5400-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 15 01:23:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 12 23:37:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	249014	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	1068352	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	534419	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	1020862	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	801361	20.00	ng/ul	0.00
86) Perylene-d12	23.49	264	856337	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.25	96	249	0.05	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.34	113	305	0.02	ng/ul	-0.06
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	10.46	131	1526	0.08	ng/ul	-0.15
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.33	176	2367	0.06	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.07	188	1020862	20.94	ng/ul	-0.10
79) Pyrene-d10	19.48	212	611	0.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.49	264	846253	19.30	ng/ul	0.14

Target Compounds	Ovalue

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

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