

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092618\
 Data File : BM016910.D
 Acq On : 26 Sep 2018 21:39
 Operator : MJ/SJ
 Sample : J5060-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 27 02:44:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 27 02:42:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	228706	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	943266	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.36	164	496402	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.10	188	1103798	20.00	ng/ul	0.00
78) Chrysene-d12	21.29	240	1201555	20.00	ng/ul	0.00
86) Perylene-d12	23.53	264	1313622	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	408	0.08	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.07	67	373	0.03	ng/ul	0.02
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	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	10.51	131	1589	0.09	ng/ul	-0.12
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	15.36	176	2948	0.08	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.20	188	3757	0.07	ng/ul	0.00
79) Pyrene-d10	19.51	212	1983	0.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	1308598	19.10	ng/ul	0.14

Target Compounds Ovalue

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

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