

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100118\
 Data File : BM016936.D
 Acq On : 02 Oct 2018 00:03
 Operator : MJ/SJ
 Sample : J5125-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B69

Quant Time: Oct 02 07:33:55 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	7.71	152	318925	20.00	ng/ul	0.00
18) Naphthalene-d8	10.49	136	1308371	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	680131	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.10	188	1349437	20.00	ng/ul	0.00
78) Chrysene-d12	21.29	240	1204444	20.00	ng/ul	-0.01
86) Perylene-d12	23.52	264	1278674	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.26	96	236	0.03	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	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	10.50	131	1622	0.07	ng/ul	-0.14
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	15.35	176	3706	0.08	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.58	200	241	0.04	ng/ul	0.11
71) Anthracene-d10	17.20	188	4026	0.06	ng/ul	0.00
79) Pyrene-d10	19.49	212	92	0.00	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.52	264	1266089	18.98	ng/ul	0.13

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

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