

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101017\
 Data File : BM012042.D
 Acq On : 10 Oct 2017 15:06
 Operator : SJ/JU
 Sample : I5726-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 11 01:13:36 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 11 01:12:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	120499	20.00	ng/ul	0.00
18) Naphthalene-d8	10.65	136	565735	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.49	164	372601	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.24	188	965093	20.00	ng/ul	0.00
75) Chrysene-d12	21.42	240	1351440	20.00	ng/ul	0.00
83) Perylene-d12	23.73	264	1603503	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	149	0.06	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	10.66	131	282	0.03	ng/ul	-0.12
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.49	176	107	0.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.24	188	965093	20.29	ng/ul	-0.09
76) Pyrene-d10	19.64	212	728	0.01	ng/ul	0.01
87) Benzo(a)pyrene-d12	23.73	264	1539307	19.96	ng/ul	0.15

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

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