

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012918\
 Data File : BM013997.D
 Acq On : 30 Jan 2018 14:58
 Operator : SJ/JU
 Sample : J1254-10
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jan 30 16:22:03 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 30 07:49:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	63226	20.00	ng/ul	0.00
18) Naphthalene-d8	10.35	136	257226	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.22	164	157774	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.98	188	371330	20.00	ng/ul	0.00
75) Chrysene-d12	21.18	240	442253	20.00	ng/ul	0.00
83) Perylene-d12	23.36	264	480337	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
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	0.00	131	0	0.00	ng/ul	
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	15.26	176	588	0.05	ng/ul	0.04
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	16.98	188	371330	20.53	ng/ul	-0.09
76) Pyrene-d10	0.00	212	0	0.00	ng/ul	
87) Benzo(a)pyrene-d12	23.36	264	470135	21.29	ng/ul	0.14

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

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