

Data Path : U:\HPCHEM1\BNA M\DATA\BM010418\
 Data File : BM013556.D
 Acq On : 04 Jan 2018 13:00
 Operator : SJ/JU
 Sample : I7103-10
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jan 04 16:02:39 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 27 03:26:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	168794	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	633355	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.30	164	349734	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	773051	20.00	ng/ul	0.00
75) Chrysene-d12	21.25	240	834316	20.00	ng/ul	0.00
83) Perylene-d12	23.47	264	838621	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	10.45	131	162	0.02	ng/ul	-0.14
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	15.30	176	495	0.02	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.06	188	773051	19.97	ng/ul	-0.11
76) Pyrene-d10	0.00	212	0	0.00	ng/ul	
87) Benzo(a)pyrene-d12	23.47	264	830268	19.44	ng/ul	0.14

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

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