

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110717\
 Data File : BM012790.D
 Acq On : 07 Nov 2017 11:46
 Operator : SJ/JU
 Sample : I6154-17
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 08 00:09:08 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 08 00:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	83229	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	349497	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	238729	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	539689	20.00	ng/ul	0.00
77) Chrysene-d12	21.37	240	490945	20.00	ng/ul	0.00
85) Perylene-d12	23.64	264	496353	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	784	0.51	ng/uL	0.00
5) Phenol-d5	6.98	99	14083	2.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	33770	10.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	42601	8.26	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	30236	5.66	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	26483	10.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	29621	10.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	55669	9.19	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	1922	0.33	ng/ul	0.01
43) Dimethylphthalate-d6	13.85	166	249745	12.64	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	240708	9.85	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	4955	1.57	ng/ul	0.02
57) Fluorene-d10	15.43	176	199435	11.79	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.57	200	32856	9.22	ng/ul	0.00
70) Anthracene-d10	17.29	188	337265	12.93	ng/ul	0.00
78) Pyrene-d10	19.58	212	333908	14.66	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.50	264	315938	13.46	ng/ul	0.00

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

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