

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101617\
 Data File : BM012224.D
 Acq On : 16 Oct 2017 20:18
 Operator : SJ/JU
 Sample : I5902-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 16 23:44:58 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 16 23:37:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	147852	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	648057	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.45	164	437370	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.20	188	1055732	20.00	ng/ul	0.00
75) Chrysene-d12	21.38	240	1014260	20.00	ng/ul	0.00
83) Perylene-d12	23.68	264	821067	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.21	96	18474	5.94	ng/uL	0.00
5) Phenol-d5	6.97	99	88410	7.37	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	187972	26.13	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	255168	25.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	185573	17.97	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	154532	31.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	189292	32.76	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	346322	31.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	18397	1.78	ng/ul	0.02
43) Dimethylphthalate-d6	13.86	166	1347393	34.86	ng/ul	0.00
46) Acenaphthylene-d8	14.15	160	1529900	32.70	ng/ul	0.00
51) 4-Nitrophenol-d4	14.69	143	34681	5.97	ng/ul	0.02
57) Fluorene-d10	15.44	176	1105810	34.85	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.59	200	213410	29.45	ng/ul	0.00
70) Anthracene-d10	17.30	188	1819948	34.86	ng/ul	0.00
76) Pyrene-d10	19.59	212	2115069	41.08	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.53	264	1450529	36.65	ng/ul	0.00

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

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