

Data Path : Z:\HPCHEM1\BNA M\DATA\BM021516\
 Data File : BM004269.D
 Acq On : 15 Feb 2016 21:05
 Operator : SJ/UM
 Sample : H1416-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9QE8

Quant Time: Feb 16 08:11:38 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 16 01:21:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	26331	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	125184	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	71462	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	142201	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	132263	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	128743	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	610	1.17	ng/uL	0.00
5) Phenol-d5	6.83	99	18356	8.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	38649	33.84	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	55514	29.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	37062	20.57	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	33278	34.57	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	38433	35.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	61461	31.67	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.72	166	205980	34.64	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	256499	35.50	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	5120	5.32	ng/ul	0.01
58) Fluorene-d10	15.30	176	174521	34.82	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	26828	30.98	ng/ul	0.00
71) Anthracene-d10	17.16	188	249670	35.91	ng/ul	0.00
79) Pyrene-d10	19.46	212	245359	34.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	242525	35.30	ng/ul	0.00

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

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