

Data Path : Z:\HPCHEM1\BNA M\Data\BM011416\
 Data File : BM003990.D
 Acq On : 13 Jan 2016 18:34
 Operator : SJ/UM
 Sample : PB87808BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jan 14 10:39:29 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 13 06:45:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	39732	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	196555	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	118015	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	265448	20.00	ng/ul	0.00
78) Chrysene-d12	21.28	240	249233	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	212939	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	4602	5.18	ng/uL	0.00
5) Phenol-d5	6.84	99	100474	30.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	60098	32.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	84980	31.34	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	87894	32.88	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	44031	29.74	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	48721	30.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	83058	28.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	132175	34.72	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	290973	31.41	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	353403	31.31	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	43384	26.07	ng/ul	0.00
58) Fluorene-d10	15.32	176	250281	31.85	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	31576	21.09	ng/ul	0.00
71) Anthracene-d10	17.17	188	393975	32.09	ng/ul	0.00
79) Pyrene-d10	19.48	212	415658	32.20	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.38	264	343487	32.30	ng/ul	0.01

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

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