

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011216\
 Data File : BM003951.D
 Acq On : 12 Jan 2016 12:50
 Operator : SJ/UM
 Sample : H1057-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G9EL0

Quant Time: Jan 12 14:00:42 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 12 00:35:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	38658	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	185045	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	112877	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	260130	20.00	ng/ul	0.00
78) Chrysene-d12	21.28	240	286630	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	258689	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3999	4.63	ng/uL	0.00
5) Phenol-d5	6.85	99	21626	6.74	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	66395	36.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	72903	27.64	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	45115	17.35	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	44842	32.17	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	48714	31.88	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	78983	29.09	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	13.73	166	313334	35.36	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	385145	35.68	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	6011	3.78	ng/ul	0.00
58) Fluorene-d10	15.32	176	281611	37.47	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	37745	25.73	ng/ul	0.00
71) Anthracene-d10	17.17	188	443129	36.83	ng/ul	0.00
79) Pyrene-d10	19.47	212	493801	33.26	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	472181	36.55	ng/ul	0.00

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

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