

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011216\
 Data File : BM003953.D
 Acq On : 12 Jan 2016 14:02
 Operator : SJ/UM
 Sample : H1057-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G9EK7

Quant Time: Jan 12 15:12:34 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	44714	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	205601	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	119863	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	256727	20.00	ng/ul	0.00
78) Chrysene-d12	21.28	240	228987	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	211986	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3640	3.64	ng/uL	0.00
5) Phenol-d5	6.85	99	22790	6.14	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	63911	30.46	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	71485	23.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	45383	15.09	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	42989	27.76	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	46038	27.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	76482	25.36	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	2474	0.62	ng/ul	0.01
44) Dimethylphthalate-d6	13.73	166	282460	30.02	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	351498	30.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.55	143	5382	3.18	ng/ul	0.00
58) Fluorene-d10	15.32	176	251560	31.52	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	31454	21.73	ng/ul	0.00
71) Anthracene-d10	17.17	188	369796	31.14	ng/ul	0.00
79) Pyrene-d10	19.47	212	381293	32.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	333179	31.47	ng/ul	0.00

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

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