

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111317\
 Data File : BM012956.D
 Acq On : 14 Nov 2017 03:52
 Operator : SJ/JU
 Sample : I6363-17
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B0FE8

Quant Time: Nov 14 04:30:27 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 14 04:22:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	106607	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	433813	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	272558	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	645007	20.00	ng/ul	0.00
77) Chrysene-d12	21.35	240	677188	20.00	ng/ul	0.00
85) Perylene-d12	23.63	264	697892	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	7559	3.83	ng/uL	0.00
5) Phenol-d5	6.96	99	162235	20.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	91636	21.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	127037	19.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	134192	19.60	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	65171	20.34	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	73676	20.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	136517	18.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	21746	3.00	ng/ul	0.01
43) Dimethylphthalate-d6	13.83	166	503353	22.32	ng/ul	0.00
46) Acenaphthylene-d8	14.11	160	569433	20.41	ng/ul	0.00
51) 4-Nitrophenol-d4	14.65	143	60587	16.76	ng/ul	0.00
57) Fluorene-d10	15.42	176	425642	22.03	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	61209	14.37	ng/ul	0.00
70) Anthracene-d10	17.27	188	694899	22.30	ng/ul	0.00
78) Pyrene-d10	19.56	212	750540	23.89	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.48	264	751388	22.77	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.98	94	25634	3.148	ng/ul 98
44) Dimethylphthalate	13.87	163	120929	5.403	ng/ul 98

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

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