

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120417\
 Data File : BM013177.D
 Acq On : 04 Dec 2017 21:22
 Operator : SJ/JU
 Sample : I6646-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0265

Quant Time: Dec 05 04:59:51 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 05 04:53:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	268970	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	1163919	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.35	164	683374	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.09	188	1536601	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1385182	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	1076767	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	11160	2.06	ng/uL	0.00
5) Phenol-d5	6.88	99	334235	14.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	202916	15.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.24	132	281050	15.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.41	113	231216	11.63	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	145846	15.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	153165	15.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.11	165	285097	14.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	263915	14.11	ng/ul	0.00
43) Dimethylphthalate-d6	13.76	166	870740	14.32	ng/ul	0.00
46) Acenaphthylene-d8	14.03	160	1139776	15.19	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	128359	12.15	ng/ul	0.00
57) Fluorene-d10	15.34	176	800015	15.30	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	96706	10.00	ng/ul	0.00
70) Anthracene-d10	17.19	188	1249556	16.24	ng/ul	0.00
76) Pyrene-d10	19.48	212	1307424	18.95	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	880712	16.06	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.90	94	39760	1.728	ng/ul#	79
44) Dimethylphthalate	13.80	163	180254	3.099	ng/ul#	97

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

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