

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM110117\
 Data File : BM012640.D
 Acq On : 01 Nov 2017 17:15
 Operator : SJ/JU
 Sample : I5937-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B0DP1

Quant Time: Nov 02 01:36:34 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 02 01:34:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	325186	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	1468327	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.47	164	1038291	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.22	188	2585081	20.00	ng/ul	0.00
75) Chrysene-d12	21.39	240	3216250	20.00	ng/ul	0.00
83) Perylene-d12	23.69	264	3637408	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	33531	5.27	ng/uL	0.00
5) Phenol-d5	7.02	99	691661	25.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.18	67	427774	26.53	ng/ul	0.00
9) 2-Chlorophenol-d4	7.38	132	537840	26.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	615236	26.50	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	302179	32.86	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	345578	36.48	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.26	165	628019	25.03	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	459289	17.36	ng/ul	0.00
43) Dimethylphthalate-d6	13.89	166	2642008	29.13	ng/ul	0.00
46) Acenaphthylene-d8	14.17	160	3025984	28.20	ng/ul	0.00
51) 4-Nitrophenol-d4	14.68	143	361243	26.79	ng/ul	0.00
57) Fluorene-d10	15.47	176	2253483	29.35	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.59	200	370161	29.59	ng/ul	0.00
70) Anthracene-d10	17.32	188	3837399	31.15	ng/ul	0.00
76) Pyrene-d10	19.61	212	4445782	30.13	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.54	264	5598980	32.40	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	7.04	94	48515	1.715	ng/ul#	84
44) Dimethylphthalate	13.93	163	555106	6.187	ng/ul	98

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

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