

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM110717\
 Data File : BM012797.D
 Acq On : 07 Nov 2017 15:55
 Operator : SJ/JU
 Sample : I6164-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET0D5

Quant Time: Nov 08 00:10:21 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 08 00:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	103447	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	445394	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	303041	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	739595	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	874349	20.00	ng/ul	0.00
85) Perylene-d12	23.64	264	848069	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	5314	2.78	ng/uL	0.00
5) Phenol-d5	6.97	99	121769	15.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	67701	16.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	104242	16.26	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	104529	15.73	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	50639	15.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	57222	15.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	112424	14.56	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	115930	15.56	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	419750	16.74	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	477684	15.40	ng/ul	0.00
51) 4-Nitrophenol-d4	14.66	143	43468	10.82	ng/ul	0.00
57) Fluorene-d10	15.43	176	357739	16.66	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.57	200	39155	8.02	ng/ul	0.00
70) Anthracene-d10	17.29	188	574615	16.08	ng/ul	0.00
78) Pyrene-d10	19.57	212	666021	16.42	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.50	264	665959	16.61	ng/ul	0.00

Target Compounds

				Ovalue
6) Phenol	7.00	94	20134	2.548 ng/ul 98
44) Dimethylphthalate	13.89	163	183063	7.356 ng/ul 100

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

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