

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120617\
 Data File : BM013233.D
 Acq On : 07 Dec 2017 05:16
 Operator : SJ/JU
 Sample : I6647-13
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0281

Quant Time: Dec 07 06:37:46 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 07 04:35:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	302691	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	1320940	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.34	164	775953	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.09	188	1756107	20.00	ng/ul	0.00
75) Chrysene-d12	21.28	240	1572849	20.00	ng/ul	0.00
83) Perylene-d12	23.51	264	1252523	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	27839	4.56	ng/uL	0.00
5) Phenol-d5	6.88	99	775939	29.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.04	67	446911	29.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.24	132	633803	30.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.41	113	664891	29.72	ng/ul	0.00
19) Nitrobenzene-d5	8.86	128	335173	32.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	367234	32.75	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.11	165	684949	29.65	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	891886	42.00	ng/ul	0.00
43) Dimethylphthalate-d6	13.76	166	2142958	31.03	ng/ul	0.00
46) Acenaphthylene-d8	14.03	160	2773342	32.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	330109	27.52	ng/ul	0.00
57) Fluorene-d10	15.34	176	1846182	31.09	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	206607	18.70	ng/ul	0.00
70) Anthracene-d10	17.19	188	2860171	32.52	ng/ul	0.00
76) Pyrene-d10	19.49	212	3011305	38.43	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.37	264	2031831	31.85	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.90	94	79887	3.086	ng/ul#	86
44) Dimethylphthalate	13.80	163	409999	6.208	ng/ul	98

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

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