

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\
 Data File : BM013111.D
 Acq On : 27 Nov 2017 14:45
 Operator : SJ/JU
 Sample : I6496-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA3

Quant Time: Nov 28 00:48:37 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 28 00:44:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	210898	20.00	ng/ul	0.00
18) Naphthalene-d8	10.49	136	935750	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.35	164	555070	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.10	188	1210896	20.00	ng/ul	0.00
75) Chrysene-d12	21.28	240	1112816	20.00	ng/ul	0.00
83) Perylene-d12	23.51	264	917539	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	16997	3.73	ng/uL	0.00
5) Phenol-d5	6.90	99	465406	25.48	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.05	67	258011	24.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	375734	25.73	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	419974	28.07	ng/ul	0.01
19) Nitrobenzene-d5	8.86	128	186544	27.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.58	143	208806	32.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	413435	25.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.63	131	451016	26.78	ng/ul	0.00
43) Dimethylphthalate-d6	13.77	166	1371371	28.17	ng/ul	0.00
46) Acenaphthylene-d8	14.04	160	1623917	26.18	ng/ul	0.00
51) 4-Nitrophenol-d4	14.58	143	221999	30.06	ng/ul	0.02
57) Fluorene-d10	15.34	176	1156627	27.95	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	72973	13.90	ng/ul	0.00
70) Anthracene-d10	17.19	188	1847475	29.80	ng/ul	0.00
76) Pyrene-d10	19.49	212	1940780	34.14	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.37	264	1481905	31.85	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.93	94	943833	52.346	ng/ul# 83
16) 4-Methylphenol	8.50	108	1757799	120.949	ng/ul 98
44) Dimethylphthalate	13.81	163	228585	4.910	ng/ul 98

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

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