

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
 Data File : BM013040.D
 Acq On : 18 Nov 2017 15:09
 Operator : SJ/JU
 Sample : I6405-09 5X
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE2T0

Quant Time: Nov 20 07:13:04 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 20 06:14:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	180774	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	697723	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.36	164	396212	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.10	188	901562	20.00	ng/ul	0.00
75) Chrysene-d12	21.29	240	1058639	20.00	ng/ul	0.00
83) Perylene-d12	23.52	264	1062996	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	1007	0.26	ng/uL	0.00
5) Phenol-d5	6.89	99	16964	1.08	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	46405	5.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	51028	4.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	30503	2.38	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	32827	6.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.58	143	32199	6.81	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	58359	4.86	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	1516	0.12	ng/ul	0.01
43) Dimethylphthalate-d6	13.77	166	197975	5.70	ng/ul	0.00
46) Acenaphthylene-d8	14.04	160	234199	5.29	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	6051	1.15	ng/ul	0.01
57) Fluorene-d10	15.35	176	162699	5.51	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	19835	5.07	ng/ul	-0.01
70) Anthracene-d10	17.20	188	255182	5.53	ng/ul	0.00
76) Pyrene-d10	19.49	212	295128	5.46	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.37	264	289789	5.38	ng/ul	0.00

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.29	88	6311	1.472	ng/uL#	79
8) Bis(2-Chloroethyl)ether	7.15	93	62584	5.378	ng/ul	96
64) N-Nitrosodiphenylamine	15.62	169	45692	1.614	ng/ul	91

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

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