

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
 Data File : BM010685.D
 Acq On : 23 Jun 2017 11:04
 Operator : SJ/JU
 Sample : I3767-01DL 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDC63DL

Quant Time: Jun 24 03:36:57 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 24 03:35:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	60048	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	272260	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.11	164	190683	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.87	188	488924	20.00	ng/ul	0.00
75) Chrysene-d12	21.08	240	544725	20.00	ng/ul	0.00
83) Perylene-d12	23.22	264	381832	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	731	0.70	ng/uL	0.01
5) Phenol-d5	6.65	99	21339	3.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.81	67	45172	13.83	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	50734	12.64	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	40219	8.48	ng/ul	0.00
19) Nitrobenzene-d5	8.60	128	34152	17.59	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	36477	16.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	72498	15.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	3770	0.76	ng/ul	0.00
43) Dimethylphthalate-d6	13.53	166	284349	16.99	ng/ul	0.00
46) Acenaphthylene-d8	13.80	160	320484	16.46	ng/ul	0.00
51) 4-Nitrophenol-d4	14.32	143	9130	3.10	ng/ul	0.00
57) Fluorene-d10	15.11	176	245599	16.97	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.23	200	44244	14.73	ng/ul	0.00
70) Anthracene-d10	16.96	188	396884	16.81	ng/ul	0.00
76) Pyrene-d10	19.27	212	461905	18.25	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.08	264	310427	16.82	ng/ul	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.10	88	23944	20.453	ng/uL# 49

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

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