

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080516\
 Data File : BG023480.D
 Acq On : 5 Aug 2016 9:50
 Operator : UM/SJ
 Sample : H4289-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Aug 05 23:14:20 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 02:24:23 2004
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	120963	20.00	ng/ul	0.00
18) Naphthalene-d8	10.96	136	545013	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	396743	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.55	188	1014822	20.00	ng/ul	0.00
75) Chrysene-d12	21.86	240	1025214	20.00	ng/ul	0.00
83) Perylene-d12	25.24	264	970819	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.52	96	154	0.05	ng/uL	0.04
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.46	67	147	0.02	ng/ul	0.02
9) 2-Chlorophenol-d4	7.56	132	57	0.01	ng/ul	-0.08
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	9.27	128	209	0.05	ng/ul	-0.02
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	10.75	165	153	0.02	ng/ul	0.18
29) 4-Chloroaniline-d4	11.19	131	126	0.01	ng/ul	0.11
43) Dimethylphthalate-d6	14.04	166	129	0.00	ng/ul	-0.13
46) Acenaphthylene-d8	14.48	160	194	0.01	ng/ul	0.00
51) 4-Nitrophenol-d4	15.00	143	172	0.03	ng/ul	0.00
57) Fluorene-d10	15.77	176	105	0.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.65	188	2740	0.06	ng/ul	0.00
76) Pyrene-d10	19.88	212	214	0.00	ng/ul	-0.05
87) Benzo(a)pyrene-d12	25.24	264	958322	21.81	ng/ul	0.24

Target Compounds	Ovalue

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

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