

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG031816\
 Data File : BG021444.D
 Acq On : 18 Mar 2016 10:53
 Operator : UM/SJ
 Sample : H1785-12DL 25X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AD6DL

Quant Time: Mar 19 00:24:07 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Mar 19 00:12:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	7315	20.00	ng/ul	0.00
18) Naphthalene-d8	10.89	136	33564	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	22280	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.44	188	58672	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	68813	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	67870	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.41	67	264	0.64	ng/ul	0.00
9) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	9.25	128	121	0.48	ng/ul	0.00
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	10.59	165	487	0.95	ng/ul	0.02
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.10	166	2582	1.39	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	2907	1.27	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.69	176	2381	1.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.54	188	4578	1.59	ng/ul	0.00
79) Pyrene-d10	19.81	212	4778	1.49	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	5060	1.62	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
20) Nitrobenzene	9.29	77	33620	47.27	ng/ul	97
73) 1,2,3,4-Tetrachlorobenzene	13.51	216	1796	2.19	ng/uL#	66

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

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