

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031716\
 Data File : BG021434.D
 Acq On : 17 Mar 2016 16:04
 Operator : UM/SJ
 Sample : H1785-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AH5

Quant Time: Mar 18 04:04:08 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 18 04:03:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	7872	20.00	ng/ul	0.00
18) Naphthalene-d8	10.89	136	37904	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.70	164	26239	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.44	188	64868	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	75214	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	73750	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.34	99	3606	4.85	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.41	67	11304	25.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.65	132	11935	21.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	8543	13.40	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	8401	29.54	ng/ul	0.00
22) 2-Nitrophenol-d4	9.98	143	9323	28.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.58	165	15568	26.79	ng/ul	0.00
29) 4-Chloroaniline-d4	11.04	131	987	1.30	ng/ul	0.00
44) Dimethylphthalate-d6	14.10	166	69049	31.60	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	79798	29.63	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	1518	4.35	ng/ul	0.02
58) Fluorene-d10	15.69	176	63293	32.13	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.81	200	12979	30.29	ng/ul	0.00
71) Anthracene-d10	17.54	188	114829	36.12	ng/ul	0.00
79) Pyrene-d10	19.81	212	134746	38.42	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	142250	41.79	ng/ul	0.00

Target Compounds Ovalue

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

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