

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100815\
 Data File : BG019070.D
 Acq On : 8 Oct 2015 17:34
 Operator : UM/NP
 Sample : PB86002BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK02

Quant Time: Oct 09 02:24:07 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 09 02:23:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	18706	20.00	ng/ul	0.00
18) Naphthalene-d8	10.84	136	80886	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	55789	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	143246	20.00	ng/ul	0.00
78) Chrysene-d12	21.69	240	175817	20.00	ng/ul	0.00
86) Perylene-d12	24.92	264	159593	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	1864	4.50	ng/uL	0.00
5) Phenol-d5	7.20	99	40888	25.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.36	67	28823	27.63	ng/ul	0.00
9) 2-Chlorophenol-d4	7.57	132	32615	26.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	34083	26.90	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	18322	29.96	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	19959	31.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	33954	26.97	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	7927	5.16	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	150891	34.30	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	155848	28.64	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	26590	29.92	ng/ul	0.00
58) Fluorene-d10	15.66	176	122048	30.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.77	200	24281	31.25	ng/ul	0.00
71) Anthracene-d10	17.51	188	214381	31.78	ng/ul	0.00
79) Pyrene-d10	19.80	212	279662	34.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.70	264	297320	36.41	ng/ul	0.00

Target Compounds	Ovalue

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

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