

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100915\
 Data File : BG019082.D
 Acq On : 9 Oct 2015 16:46
 Operator : UM/NP
 Sample : PB86051BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Oct 10 00:38:50 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 10 00:37:29 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	19515	20.00	ng/ul	0.00
18) Naphthalene-d8	10.84	136	90587	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	64840	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	173298	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	216904	20.00	ng/ul	0.00
86) Perylene-d12	24.92	264	201879	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	2335	5.41	ng/uL	0.00
5) Phenol-d5	7.19	99	49671	29.54	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.36	67	34104	31.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.57	132	38795	30.73	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	43051	32.56	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	21583	31.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	24219	34.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	42451	30.10	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	1718	1.00	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	187768	36.73	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	197634	31.25	ng/ul	0.00
52) 4-Nitrophenol-d4	14.85	143	33790	32.72	ng/ul	0.00
58) Fluorene-d10	15.65	176	155393	33.57	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.76	200	29772	31.67	ng/ul	0.00
71) Anthracene-d10	17.51	188	272333	33.37	ng/ul	0.00
79) Pyrene-d10	19.80	212	342220	34.06	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.70	264	351268	34.01	ng/ul	0.00

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

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