

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121016\
 Data File : BG025086.D
 Acq On : 11 Dec 2016 2:27
 Operator : UM/SJ
 Sample : H5861-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Dec 12 04:27:00 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 04:23:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	84031	20.00	ng/ul	0.00
18) Naphthalene-d8	11.06	136	370698	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.85	164	306346	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.60	188	714157	20.00	ng/ul	0.00
75) Chrysene-d12	21.89	240	889495	20.00	ng/ul	0.00
83) Perylene-d12	25.30	264	900546	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.59	96	2418	1.49	ng/uL	0.00
5) Phenol-d5	7.37	99	54498	7.52	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.55	67	149648	33.37	ng/ul	0.00
9) 2-Chlorophenol-d4	7.76	132	132172	26.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.94	113	106017	17.50	ng/ul	0.00
19) Nitrobenzene-d5	9.41	128	87592	33.37	ng/ul	0.00
22) 2-Nitrophenol-d4	10.14	143	107507	32.85	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.68	165	191099	30.36	ng/ul	0.00
29) 4-Chloroaniline-d4	11.21	131	2894	0.47	ng/ul	0.01
43) Dimethylphthalate-d6	14.25	166	718910	34.12	ng/ul	0.00
46) Acenaphthylene-d8	14.55	160	800964	34.70	ng/ul	0.00
51) 4-Nitrophenol-d4	15.05	143	27726	7.66	ng/ul	0.00
57) Fluorene-d10	15.84	176	673897	33.98	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.96	200	152788	34.60	ng/ul	0.00
70) Anthracene-d10	17.70	188	1082889	35.93	ng/ul	0.00
76) Pyrene-d10	19.97	212	1340881	36.60	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.07	264	1344282	36.84	ng/ul	0.00

Target Compounds Ovalue

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

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