

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030816\
 Data File : BG021374.D
 Acq On : 8 Mar 2016 19:51
 Operator : UM/SJ
 Sample : H1705-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MS-GW-E350

Quant Time: Mar 09 01:05:14 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 09 01:04:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	9568	20.00	ng/ul	0.00
18) Naphthalene-d8	10.93	136	46463	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	28015	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	62797	20.00	ng/ul	0.00
78) Chrysene-d12	21.73	240	67656	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	63741	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	255	1.11	ng/uL	0.00
5) Phenol-d5	7.27	99	7426	7.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.44	67	22299	33.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	19351	27.65	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	14440	18.24	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	12626	33.97	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	13548	34.34	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	22826	32.27	ng/ul	0.00
29) 4-Chloroaniline-d4	11.06	131	1475	1.67	ng/ul	0.00
44) Dimethylphthalate-d6	14.13	166	85462	36.90	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	108452	37.10	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	2904	6.37	ng/ul	0.00
58) Fluorene-d10	15.72	176	76000	37.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	13734	31.16	ng/ul	0.00
71) Anthracene-d10	17.57	188	119605	38.64	ng/ul	0.00
79) Pyrene-d10	19.84	212	133357	39.43	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.76	264	125135	36.89	ng/ul	0.00

Target Compounds Qvalue

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

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