

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030416\
 Data File : BG021295.D
 Acq On : 5 Mar 2016 7:27
 Operator : UM/SJ
 Sample : H1650-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MS-RB-01

Quant Time: Mar 07 00:51:16 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Mar 05 05:52:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	10169	20.00	ng/ul	0.00
18) Naphthalene-d8	10.93	136	50039	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	29083	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	69316	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	76772	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	69656	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.53	96	207	0.83	ng/uL	0.00
5) Phenol-d5	7.27	99	6474	6.01	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	23847	33.20	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	19170	25.51	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	13222	15.55	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	13348	32.93	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	14145	32.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	22349	28.94	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	848	0.87	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	85626	35.13	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	110285	36.36	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	2553	5.37	ng/ul	0.00
58) Fluorene-d10	15.72	176	78931	36.88	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	14612	29.37	ng/ul	0.00
71) Anthracene-d10	17.57	188	125312	36.35	ng/ul	0.00
79) Pyrene-d10	19.85	212	141246	36.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	134864	35.86	ng/ul	0.00

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

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