

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
 Data File : BG021353.D
 Acq On : 8 Mar 2016 00:35
 Operator : UM/SJ
 Sample : H1625-16 10X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHFH3

Quant Time: Mar 08 01:26:58 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 00:45:15 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	10501	20.00	ng/ul	0.00
18) Naphthalene-d8	10.93	136	49606	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.73	164	28616	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	66344	20.00	ng/ul	0.00
78) Chrysene-d12	21.73	240	70358	20.00	ng/ul	0.00
86) Perylene-d12	24.98	264	66137	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.28	99	1606	1.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.44	67	1198	1.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	1146	1.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	1155	1.32	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	596	1.48	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	493	1.15	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	971	1.27	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	781	0.80	ng/ul	0.01
44) Dimethylphthalate-d6	14.14	166	4002	1.67	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	5350	1.79	ng/ul	0.00
52) 4-Nitrophenol-d4	14.93	143	218	0.47	ng/ul	0.01
58) Fluorene-d10	15.72	176	4240	2.01	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.57	188	5938	1.80	ng/ul	0.00
79) Pyrene-d10	19.84	212	6583	1.85	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.75	264	6367	1.78	ng/ul	0.00

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

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