

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032116\
 Data File : BG021455.D
 Acq On : 21 Mar 2016 22:23
 Operator : UM/SJ
 Sample : H1835-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T21

Quant Time: Mar 22 01:45:14 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 22 01:42:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	10259	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	47288	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	29227	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	69321	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	82960	20.00	ng/ul	0.00
86) Perylene-d12	24.90	264	80533	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	218	0.95	ng/uL	0.00
5) Phenol-d5	7.34	99	5504	5.67	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.40	67	14959	26.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	16067	21.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	10462	12.59	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	10531	29.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	11773	29.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	18330	25.29	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.09	166	74853	30.76	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	88525	29.51	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	1728	4.44	ng/ul	0.02
58) Fluorene-d10	15.68	176	68875	31.39	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	12901	28.17	ng/ul	0.00
71) Anthracene-d10	17.53	188	123424	36.33	ng/ul	0.00
79) Pyrene-d10	19.80	212	143103	36.99	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	150404	40.47	ng/ul	0.00

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

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