

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032116\
 Data File : BG021451.D
 Acq On : 21 Mar 2016 19:46
 Operator : UM/SJ
 Sample : H1835-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T13

Quant Time: Mar 22 01:44:36 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	8.07	152	10217	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	46341	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	27853	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	62439	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	77430	20.00	ng/ul	0.00
86) Perylene-d12	24.90	264	72786	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	7.35	99	5331	5.52	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.40	67	16059	28.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.65	132	16018	21.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	10533	12.73	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	10877	31.28	ng/ul	0.00
22) 2-Nitrophenol-d4	9.97	143	12175	30.64	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	19472	27.41	ng/ul	0.00
29) 4-Chloroaniline-d4	11.03	131	964	1.04	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	74246	32.01	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	91591	32.04	ng/ul	0.00
52) 4-Nitrophenol-d4	15.10	143	1906	5.14	ng/ul	0.03
58) Fluorene-d10	15.68	176	67921	32.48	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	12268	29.74	ng/ul	0.00
71) Anthracene-d10	17.52	188	119076	38.92	ng/ul	0.00
79) Pyrene-d10	19.80	212	144218	39.94	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	150801	44.89	ng/ul	0.00

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

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