

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032116\
 Data File : BG021456.D
 Acq On : 21 Mar 2016 23:03
 Operator : UM/SJ
 Sample : H1835-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T22

Quant Time: Mar 22 01:45:20 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	9462	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	47758	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	30444	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	69051	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	81392	20.00	ng/ul	0.00
86) Perylene-d12	24.90	264	78729	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.34	99	4581	5.12	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	13209	24.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.65	132	13682	20.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	9424	12.30	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	9252	25.82	ng/ul	0.00
22) 2-Nitrophenol-d4	9.97	143	10520	25.69	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.58	165	17260	23.57	ng/ul	0.00
29) 4-Chloroaniline-d4	11.03	131	1604	1.68	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	72926	28.77	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	81588	26.11	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	1806	4.46	ng/ul	0.02
58) Fluorene-d10	15.68	176	62948	27.54	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	12085	26.49	ng/ul	0.00
71) Anthracene-d10	17.52	188	114843	33.94	ng/ul	0.00
79) Pyrene-d10	19.80	212	138828	36.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	143739	39.56	ng/ul	0.00

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

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