

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032116\
 Data File : BG021457.D
 Acq On : 21 Mar 2016 23:42
 Operator : UM/SJ
 Sample : H1835-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T23

Quant Time: Mar 22 01:45:26 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	10130	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	50454	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	31489	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	76217	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	90988	20.00	ng/ul	-0.01
86) Perylene-d12	24.89	264	89039	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	4929	5.15	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.40	67	14406	25.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	15059	20.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	10247	12.49	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	10389	27.45	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	11841	27.37	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.58	165	18800	24.31	ng/ul	0.00
29) 4-Chloroaniline-d4	11.03	131	1372	1.36	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	84400	32.19	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	97144	30.06	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	1631	3.89	ng/ul	0.03
58) Fluorene-d10	15.68	176	73517	31.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	12940	25.70	ng/ul	0.00
71) Anthracene-d10	17.52	188	123213	32.99	ng/ul	0.00
79) Pyrene-d10	19.79	212	147443	34.75	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	153142	37.27	ng/ul	0.00

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

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