

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021915\
 Data File : BG016013.D
 Acq On : 19 Feb 2015 21:41
 Operator : TP/IZ
 Sample : MDL-06-W
 Misc : MDL-WATER-3PPM-SOM0
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Feb 20 06:35:39 2015

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-MA-BG021915.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Feb 20 05:38:39 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	93789	20.00	ng/ul	0.00
16) Naphthalene-d8	10.60	136	409698	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.44	164	246727	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.18	188	532643	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	571604	20.00	ng/ul	0.00
83) Perylene-d12	23.65	264	561694	20.00	ng/ul	0.00

System Monitoring Compounds

3) Phenol-d5	6.97	99	286627	32.12	ng/ul	0.00
5) Bis-(2-Chloroethyl)ether-d	7.14	67	177644	33.71	ng/ul	0.00
7) 2-Chlorophenol-d4	7.34	132	222068	33.64	ng/ul	0.00
11) 4-Methylphenol-d8	8.51	113	216417	32.77	ng/ul	0.00
17) Nitrobenzene-d5	8.96	128	108818	37.58	ng/ul	0.00
20) 2-Nitrophenol-d4	9.69	143	98774	36.98	ng/ul	0.00
24) 2,4-Dichlorophenol-d3	10.22	165	184196	31.79	ng/ul	0.00
27) 4-Chloroaniline-d4	10.73	131	338245	52.99	ng/ul	0.00
41) Dimethylphthalate-d6	13.85	166	620362	37.61	ng/ul	0.00
44) Acenaphthylene-d8	14.13	160	801209	34.32	ng/ul	0.00
49) 4-Nitrophenol-d4	14.62	143	101031	30.46	ng/ul	0.00
55) Fluorene-d10	15.43	176	587607	35.44	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	15.54	200	40403	19.67	ng/ul	0.00
68) Anthracene-d10	17.28	188	900337	35.64	ng/ul	0.00
76) Pyrene-d10	19.57	212	976413	36.37	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.50	264	923339	36.25	ng/ul	0.00

Target Compounds

					Qvalue
70) 1,2,3,4-Tetrachlorobenzene	13.23	216	15011	2.22 ng/uL	94
71) Pentachlorobenzene	14.75	250	30718	4.63 ng/uL	90

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

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