

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021915\
 Data File : BG016007.D
 Acq On : 19 Feb 2015 18:14
 Operator : TP/IZ
 Sample : MDL-BLK-W
 Misc : MDL-WATER-BLK-SOM0
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Feb 20 09:31:10 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	7.82	152	98905	20.00	ng/ul	0.00
16) Naphthalene-d8	10.60	136	435539	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.44	164	263125	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.18	188	560029	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	593008	20.00	ng/ul	0.00
83) Perylene-d12	23.65	264	589346	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) Phenol-d5	6.97	99	315465	33.53	ng/ul	0.00
5) Bis-(2-Chloroethyl)ether-d	7.15	67	193742	34.86	ng/ul	0.00
7) 2-Chlorophenol-d4	7.35	132	241652	34.71	ng/ul	0.00
11) 4-Methylphenol-d8	8.51	113	241472	34.67	ng/ul	0.00
17) Nitrobenzene-d5	8.97	128	123838	40.23	ng/ul	0.00
20) 2-Nitrophenol-d4	9.68	143	113199	39.87	ng/ul	0.00
24) 2,4-Dichlorophenol-d3	10.21	165	207114	33.62	ng/ul	0.00
27) 4-Chloroaniline-d4	10.74	131	371938	54.81	ng/ul	0.00
41) Dimethylphthalate-d6	13.85	166	673630	38.29	ng/ul	0.00
44) Acenaphthylene-d8	14.13	160	875213	35.15	ng/ul	0.00
49) 4-Nitrophenol-d4	14.62	143	120027	33.93	ng/ul	0.00
55) Fluorene-d10	15.43	176	631933	35.74	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	15.54	200	51425	23.82	ng/ul	0.00
68) Anthracene-d10	17.27	188	959459	36.12	ng/ul	0.00
76) Pyrene-d10	19.57	212	1025152	36.80	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.51	264	968791	36.25	ng/ul	0.00

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

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