

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120114\
 Data File : BG015558.D
 Acq On : 1 Dec 2014 12:42
 Operator : TP/JJ
 Sample : F4638-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBWY2

Quant Time: Dec 02 01:02:04 2014
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG112514.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 02 00:56:15 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	103022	20.00	ng/ul	0.00
16) Naphthalene-d8	10.55	136	407048	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.40	164	298508	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.15	188	700309	20.00	ng/ul	0.00
73) Chrysene-d12	21.33	240	982757	20.00	ng/ul	0.00
81) Perylene-d12	23.62	264	962269	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) Phenol-d5	0.00	99	0	0.00	ng/ul	
5) Bis-(2-Chloroethyl)ether-d	7.11	67	179	0.03	ng/ul	0.00
7) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
11) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
17) Nitrobenzene-d5	8.91	128	97	0.03	ng/ul	0.00
20) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
24) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
27) 4-Chloroaniline-d4	10.57	131	401	0.05	ng/ul	-0.12
41) Dimethylphthalate-d6	13.80	166	384	0.02	ng/ul	-0.01
44) Acenaphthylene-d8	13.97	160	267	0.01	ng/ul	-0.13
49) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
55) Fluorene-d10	15.40	176	2165	0.11	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
68) Anthracene-d10	17.25	188	2156	0.07	ng/ul	0.00
74) Pyrene-d10	19.55	212	548	0.01	ng/ul	0.00
85) Benzo(a)pyrene-d12	23.47	264	462	0.01	ng/ul	0.00

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

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