

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120114\
 Data File : BG015561.D
 Acq On : 1 Dec 2014 14:26
 Operator : TP/JJ
 Sample : F4757-13DL2 100X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 F3W22DL2

Quant Time: Dec 02 01:05:22 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	88293	20.00	ng/ul	0.00
16) Naphthalene-d8	10.56	136	363328	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.40	164	270060	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.15	188	652049	20.00	ng/ul	0.00
73) Chrysene-d12	21.33	240	852305	20.00	ng/ul	0.00
81) Perylene-d12	23.61	264	795887	20.00	ng/ul	-0.01

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	181	0.03	ng/ul	0.00
7) 2-Chlorophenol-d4	7.33	132	385	0.07	ng/ul	0.03
11) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
17) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
20) 2-Nitrophenol-d4	9.46	143	149	0.05	ng/ul	-0.17
24) 2,4-Dichlorophenol-d3	10.21	165	144	0.03	ng/ul	0.04
27) 4-Chloroaniline-d4	10.56	131	749	0.10	ng/ul	-0.13
41) Dimethylphthalate-d6	13.68	166	148	0.01	ng/ul	-0.14
44) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
49) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
55) Fluorene-d10	15.39	176	1905	0.11	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
68) Anthracene-d10	17.24	188	898	0.03	ng/ul	-0.01
74) Pyrene-d10	19.43	212	74	0.00	ng/ul	-0.11
85) Benzo(a)pyrene-d12	23.48	264	397	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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