

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062215\
 Data File : BG017742.D
 Acq On : 22 Jun 2015 12:34
 Operator : TP/IZ
 Sample : G2653-05
 Misc : LOD-SVOC-WATER-2PPM
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-WATER2015-01

Quant Time: Jun 23 03:31:53 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 06:05:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	65136	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	285675	20.00	ng	-0.01
38) Acenaphthene-d10	14.26	164	176033	20.00	ng	-0.01
63) Phenanthrene-d10	17.02	188	401444	20.00	ng	-0.01
75) Chrysene-d12	21.22	240	430320	20.00	ng	-0.01
86) Perylene-d12	23.45	264	399948	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	466213	125.61	ng	0.00
7) Phenol-d6	6.79	99	627672	120.32	ng	-0.01
23) Nitrobenzene-d5	8.76	82	383369	77.19	ng	-0.01
41) 2,4,6-Tribromophenol	15.76	330	169903	104.09	ng	-0.01
44) 2-Fluorobiphenyl	12.88	172	774702	71.48	ng	-0.02
78) Terphenyl-d14	19.67	244	1319691	69.69	ng	-0.01
Target Compounds						Qvalue
6) Aniline	6.95	93	5651	0.79	ng	# 86
10) Phenol	6.82	94	8088	1.42	ng	94
33) 4-Chloroaniline	10.55	127	5502	0.81	ng	# 87
40) Hexachlorocyclopentadiene	12.42	237	3004	0.95	ng	93
53) 2,4-Dinitrophenol	14.38	184	841	7.77	ng	# 84
64) 4,6-Dinitro-2-methylphenol	15.40	198	1405	7.73	ng	# 73

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

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