

Data Path : S:\HPCHEM1\BNA_G\DATA\BG010615\
 Data File : BG015857.D
 Acq On : 6 Jan 2015 12:57
 Operator : TP/IZ
 Sample : G1001-05
 Misc : LOD-WATER-5PPM
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-WATER2015-01

Quant Time: Jan 07 17:45:23 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG010515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 06 08:46:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	25468	20.00	ng	0.00
21) Naphthalene-d8	10.51	136	96935	20.00	ng	0.00
38) Acenaphthene-d10	14.37	164	72254	20.00	ng	0.00
63) Phenanthrene-d10	17.11	188	177409	20.00	ng	0.00
75) Chrysene-d12	21.29	240	274461	20.00	ng	0.00
86) Perylene-d12	23.55	264	257353	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	170664	57.63	ng	0.00
7) Phenol-d6	6.91	99	232076	56.27	ng	0.00
23) Nitrobenzene-d5	8.88	82	164381	33.27	ng	0.00
41) 2,4,6-Tribromophenol	15.86	330	145420	52.95	ng	0.00
44) 2-Fluorobiphenyl	12.98	172	352566	34.85	ng	0.00
78) Terphenyl-d14	19.74	244	855617	34.83	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.25	88	1838	2.71	ng	# 41
3) Pyridine	3.65	79	5686	3.05	ng	89
4) n-Nitrosodimethylamine	3.57	42	2635	3.29	ng	# 69
40) Hexachlorocyclopentadiene	12.53	237	3675	4.72	ng	94
55) 4-Nitrophenol	14.59	139	2597	2.88	ng	# 55
64) 4,6-Dinitro-2-methylphenol	15.49	198	1630	4.70	ng	85
69) Pentachlorophenol	16.77	266	2483	4.44	ng	# 58

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

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