

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100317\
 Data File : BG029006.D
 Acq On : 3 Oct 2017 15:27
 Operator : SJ/JU
 Sample : I5275-05
 Misc : LOD 5 PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-WATER-05-QT4-2017

Quant Time: Oct 05 01:20:09 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 16:57:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.24	152	33838	20.00	ng	-0.09
21) Naphthalene-d8	11.06	136	140803	20.00	ng	-0.09
38) Acenaphthene-d10	14.87	164	102923	20.00	ng	-0.08
63) Phenanthrene-d10	17.62	188	259891	20.00	ng	-0.08
75) Chrysene-d12	21.92	240	285949	20.00	ng	-0.11
86) Perylene-d12	25.34	264	276756	20.00	ng	-0.19
System Monitoring Compounds						
5) 2-Fluorophenol	5.81	112	345571	168.51	ng	-0.10
7) Phenol-d6	7.40	99	463806	150.21	ng	-0.09
23) Nitrobenzene-d5	9.41	82	252672	85.84	ng	-0.10
41) 2,4,6-Tribromophenol	16.36	330	287390	168.83	ng	-0.08
44) 2-Fluorobiphenyl	13.48	172	663222	85.92	ng	-0.08
78) Terphenyl-d14	20.20	244	1114421	83.11	ng	-0.08
Target Compounds						
3) Pyridine	4.13	79	11718	4.947	ng	86
4) n-Nitrosodimethylamine	4.04	42	6001	5.569	ng	# 71
32) Benzoic acid	10.36	122	1221	6.678	ng	# 49
33) 4-Chloroaniline	11.23	127	15487	5.027	ng	# 83
35) Caprolactam	11.99	113	4318	4.773	ng	# 81
52) 3-Nitroaniline	14.78	138	8737	5.145	ng	# 84
55) 4-Nitrophenol	15.14	139	3356	2.697	ng	# 90
64) 4,6-Dinitro-2-methylphenol	16.01	198	5323	3.174	ng	87

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

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