

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\
 Data File : BG032571.D
 Acq On : 6 Feb 2018 18:25
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD-W-8 ppm
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-WATER-01-QT1-2018

Quant Time: Feb 07 13:56:02 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 18:57:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.71	152	21696	20.00	ng	0.00
21) Naphthalene-d8	11.58	136	86054	20.00	ng	0.00
38) Acenaphthene-d10	15.34	164	63061	20.00	ng	0.00
63) Phenanthrene-d10	18.07	188	158242	20.00	ng	0.00
75) Chrysene-d12	22.40	240	208133	20.00	ng	-0.01
86) Perylene-d12	26.07	264	233341	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.14	112	119144	110.30	ng	0.00
7) Phenol-d6	7.82	99	149811	103.93	ng	0.00
23) Nitrobenzene-d5	9.90	82	92458	67.11	ng	0.00
41) 2,4,6-Tribromophenol	16.82	330	134089	111.71	ng	0.00
44) 2-Fluorobiphenyl	13.96	172	288543	54.38	ng	0.00
78) Terphenyl-d14	20.61	244	765449	78.78	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	13.53	237	5582	3.028	ng	85
53) 2,4-Dinitrophenol	15.45	184	355	4.546	ng	# 58
69) Pentachlorophenol	17.71	266	4092	2.518	ng	# 79

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

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