

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072418\
 Data File : BG035855.D
 Acq On : 24 Jul 2018 19:57
 Operator : JU/SJ
 Sample : J3762-07
 Misc : LOD 8PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-WATER-01-QT3-2018

Quant Time: Jul 25 03:16:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 24 14:48:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	39398	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	141090	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	97626	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	262716	20.00	ng	0.00
75) Chrysene-d12	21.66	240	295700	20.00	ng	0.00
86) Perylene-d12	24.85	264	281919	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	297836	125.51	ng	0.00
7) Phenol-d6	7.19	99	420492	118.76	ng	0.00
23) Nitrobenzene-d5	9.19	82	282796	83.73	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	175159	136.12	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	641868	88.08	ng	0.00
78) Terphenyl-d14	20.00	244	1375990	102.57	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.83	237	9994	5.172	ng	91
53) 2,4-Dinitrophenol	14.81	184	88	6.672	ng	# 1
69) Pentachlorophenol	17.06	266	3256	1.703	ng	# 63

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

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