Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072418\
Data File : BG035864.D
Acq On : 25 Jul 2018 2:27
Operator : JU/SJ
Sample : J3762-09
Misc : MDL 8PPM
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MDL-WATER-03-QT3-2018

Quant Time: Jul 25 04:14:04 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	35891	20.00	ng		0.00
21) Naphthalene-d8	10.83	136	136197	20.00	ng		0.00
38) Acenaphthene-d10	14.65	164	95941	20.00	ng		0.00
63) Phenanthrene-d10	17.39	188	256855	20.00	ng		0.00
75) Chrysene-d12	21.66	240	298456	20.00	ng		0.00
86) Perylene-d12	24.85	264	278053	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.61	112	281156	130.06	ng		0.00
7) Phenol-d6	7.19	99	400152	124.06	ng		0.00
23) Nitrobenzene-d5	9.19	82	270667	83.02	ng		0.00
41) 2,4,6-Tribromophenol	16.13	330	170237	134.62	ng		0.00
44) 2-Fluorobiphenyl	13.27	172	639533	89.31	ng		0.00
78) Terphenyl-d14	20.00	244	1375382	101.58	ng		0.00
Target Compounds							Qvalue
40) Hexachlorocyclopentadiene	12.83	237	6785	3.573	ng		86
53) 2,4-Dinitrophenol	14.36	184	58	6.643	ng	#	1
69) Pentachlorophenol	17.06	266	1502	0.804	ng	#	70

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

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