Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
Data File : BG037460.D
Acq On : 19 Oct 2018 19:12
Operator : JU/SJ
Sample : J5164-07
Misc : LOD 8PPM
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : LOD-MDL-WATER-01-QT4-2018

Quant Time: Oct 20 02:37:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	46190	20.00	ng		0.00
21) Naphthalene-d8	10.93	136	210437	20.00	ng		0.00
39) Acenaphthene-d10	14.74	164	137389	20.00	ng		0.00
64) Phenanthrene-d10	17.49	188	332132	20.00	ng		0.00
76) Chrysene-d12	21.77	240	311768	20.00	ng		0.00
87) Perylene-d12	25.11	264	310996	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.65	112	448331	162.27	ng		0.00
7) Phenol-d6	7.25	99	654797	156.74	ng		0.00
23) Nitrobenzene-d5	9.28	82	363339	96.72	ng		0.00
42) 2,4,6-Tribromophenol	16.23	330	245584	163.89	ng		0.00
45) 2-Fluorobiphenyl	13.37	172	874847	96.02	ng		0.00
79) Terphenyl-d14	20.09	244	1356182	92.95	ng		0.00
Target Compounds							Qvalue
41) Hexachlorocyclopentadiene	12.90	237	20137	9.513	ng		97
54) 2,4-Dinitrophenol	14.85	184	1885	9.503	ng	#	90
70) Pentachlorophenol	17.13	266	11981	4.732	ng		94

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

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