Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101818\
Data File : BG037434.D
Acq On : 19 Oct 2018 00:28
Operator : JU/SJ
Sample : J5164-07
Misc : LOD 8 PPM
ALS Vial : 16 Sample Multiplier: 1

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

Quant Time: Oct 19 08:57:17 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	43770	20.00	ng		0.00
21) Naphthalene-d8	10.93	136	199452	20.00	ng		0.00
39) Acenaphthene-d10	14.74	164	132877	20.00	ng		0.00
64) Phenanthrene-d10	17.49	188	328257	20.00	ng		0.00
76) Chrysene-d12	21.77	240	308161	20.00	ng		0.00
87) Perylene-d12	25.10	264	305616	20.00	ng		-0.01
System Monitoring Compounds							
5) 2-Fluorophenol	5.65	112	434968	166.14	ng		0.00
7) Phenol-d6	7.25	99	627386	158.48	ng		0.00
23) Nitrobenzene-d5	9.28	82	347356	97.56	ng		0.00
42) 2,4,6-Tribromophenol	16.23	330	242567	167.37	ng		0.00
45) 2-Fluorobiphenyl	13.37	172	840974	95.44	ng		0.00
79) Terphenyl-d14	20.08	244	1328095	92.09	ng		0.00
Target Compounds							Qvalue
41) Hexachlorocyclopentadiene	12.90	237	19134	9.346	ng		96
54) 2,4-Dinitrophenol	14.85	184	1201	8.380	ng	#	1
70) Pentachlorophenol	17.12	266	11365	4.541	ng		89

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

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