Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102318\
Data File : BF110094.D
Acq On : 23 Oct 2018 22:54
Operator : JU/SJ
Sample : J5164-07
Misc : LOD 8PPM
ALS Vial : 16 Sample Multiplier: 1

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

Quant Time: Oct 24 07:34:59 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	187466	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	790563	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	390345	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	742895	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	634110	20.00	ng	-0.02
87) Perylene-d12	15.67	264	539976	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	1256259	109.13	ng	-0.02
7) Phenol-d6	6.59	99	1571701	106.74	ng	-0.02
23) Nitrobenzene-d5	7.53	82	960045	72.03	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	417771	108.05	ng	-0.01
45) 2-Fluorobiphenyl	9.32	172	1755681	70.63	ng	-0.02
79) Terphenyl-d14	13.09	244	1839758	60.34	ng	-0.02
Target Compounds						Qvalue
41) Hexachlorocyclopentadiene	9.11	237	43578	7.007	ng	98
54) 2,4-Dinitrophenol	10.06	184	9949	8.953	ng	90
70) Pentachlorophenol	11.30	266	24307	4.875	ng	97

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

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