Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
Data File : BM016052.D
Acq On : 24 Jul 2018 07:15
Operator : SJ/JU
Sample : J3762-09
Misc : MDL 8 PPM
ALS Vial : 22 Sample Multiplier: 1

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

Quant Time: Jul 24 09:39:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 23 16:08:38 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	39111	20.00	ng		0.00
21) Naphthalene-d8	11.05	136	152105	20.00	ng		0.00
38) Acenaphthene-d10	14.85	164	81094	20.00	ng		0.00
63) Phenanthrene-d10	17.58	188	173201	20.00	ng		0.00
75) Chrysene-d12	21.70	240	192760	20.00	ng		0.00
86) Perylene-d12	24.07	264	204406	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.74	112	305682	129.83	ng		0.00
7) Phenol-d6	7.38	99	381357	124.03	ng		0.00
23) Nitrobenzene-d5	9.41	82	293686	89.80	ng		0.00
41) 2,4,6-Tribromophenol	16.33	330	133568	129.86	ng		0.00
44) 2-Fluorobiphenyl	13.47	172	594342	89.18	ng		0.00
78) Terphenyl-d14	20.15	244	1132921	123.03	ng		0.00
Target Compounds							Qvalue
40) Hexachlorocyclopentadiene	13.01	237	4047	10.372	ng		87
53) 2,4-Dinitrophenol	15.04	184	1308	8.089	ng	#	57
69) Pentachlorophenol	17.24	266	3907	2.921	ng		94

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

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