

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052218\
 Data File : BM015255.D
 Acq On : 23 May 2018 00:16
 Operator : SJ/JU
 Sample : J2133-07
 Misc : LOD 8 PPM
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT2-2018

Quant Time: May 23 05:39:47 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM052218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 22 15:47:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.40	152	200921	20.00	ng	0.00
21) Naphthalene-d8	11.24	136	817723	20.00	ng	0.00
38) Acenaphthene-d10	15.02	164	422500	20.00	ng	0.00
63) Phenanthrene-d10	17.73	188	872499	20.00	ng	0.00
75) Chrysene-d12	21.83	240	881073	20.00	ng	0.00
86) Perylene-d12	24.27	264	888928	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.89	112	1653039	134.61	ng	0.00
7) Phenol-d6	7.55	99	2113758	132.18	ng	0.00
23) Nitrobenzene-d5	9.59	82	1302815	87.28	ng	0.00
41) 2,4,6-Tribromophenol	16.49	330	745893	140.79	ng	0.00
44) 2-Fluorobiphenyl	13.65	172	2872792	84.53	ng	0.00
78) Terphenyl-d14	20.29	244	3770180	94.49	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	13.20	237	39761	9.090	ng	99
53) 2,4-Dinitrophenol	15.13	184	15023	9.827	ng	90
69) Pentachlorophenol	17.39	266	41592	6.656	ng	98

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

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