

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020718\
 Data File : BM014179.D
 Acq On : 08 Feb 2018 02:05
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD 8 PPM
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT1-2018

Quant Time: Feb 08 08:21:31 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 17:16:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	79679	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	372799	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	254420	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	662722	20.00	ng	0.00
75) Chrysene-d12	21.16	240	771303	20.00	ng	0.00
86) Perylene-d12	23.34	264	704250	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	657108	143.74	ng	0.00
7) Phenol-d6	6.73	99	900262	139.05	ng	0.00
23) Nitrobenzene-d5	8.70	82	872901	104.59	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	534127	147.95	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	2099768	103.02	ng	0.00
78) Terphenyl-d14	19.60	244	3903433	98.89	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.33	237	22055	10.07	ng	95
53) 2,4-Dinitrophenol	14.37	184	8769	10.65	ng	89
69) Pentachlorophenol	16.62	266	27601	10.31	ng	88

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

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