

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022318\
 Data File : BF103197.D
 Acq On : 23 Feb 2018 14:11
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD-S-8ppm
 ALS Vial : 15 Sample Multiplier: 1

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

Quant Time: Feb 23 17:05:04 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	172849	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	748533	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	351197	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	605290	20.00	ng	0.00
75) Chrysene-d12	14.04	240	459789	20.00	ng	0.00
86) Perylene-d12	15.54	264	438297	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1260659	110.31	ng	0.00
7) Phenol-d6	6.50	99	1557707	111.39	ng	0.00
23) Nitrobenzene-d5	7.44	82	1149484	87.70	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	343257	115.47	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1720907	84.21	ng	0.00
78) Terphenyl-d14	12.99	244	1433304	74.94	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	9.02	237	7400	10.826	ng	93
53) 2,4-Dinitrophenol	9.98	184	8873	13.276	ng	# 19
69) Pentachlorophenol	11.20	266	20440	6.177	ng	100

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

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