

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020718\
 Data File : BM014176.D
 Acq On : 08 Feb 2018 00:18
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD 1 PPM
 ALS Vial : 12 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:50:40 PM

Quant Time: Feb 08 11:45:39 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	81618	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	346751	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	232307	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	554540	20.00	ng	0.00
75) Chrysene-d12	21.16	240	600406	20.00	ng	0.00
86) Perylene-d12	23.34	264	545656	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	612281	130.75	ng	0.00
7) Phenol-d6	6.73	99	824687	124.35	ng	0.00
23) Nitrobenzene-d5	8.70	82	711252	91.62	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	422194	128.08	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	1633672	87.79	ng	0.00
78) Terphenyl-d14	19.60	244	2659252	86.54	ng	0.00

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

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