

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
 Data File : BM014188.D
 Acq On : 08 Feb 2018 11:26
 Operator : SJ/JU
 Sample : J1161-09
 Misc : LOD 8 PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-03-QT1-2018

Quant Time: Feb 08 17:12:32 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	7.55	152	67172	20.00	ng	0.00
21) Naphthalene-d8	10.31	136	323750	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	219474	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	566724	20.00	ng	0.00
75) Chrysene-d12	21.16	240	652897	20.00	ng	0.00
86) Perylene-d12	23.34	264	580271	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	566001	146.87	ng	0.00
7) Phenol-d6	6.73	99	776855	142.33	ng	0.00
23) Nitrobenzene-d5	8.70	82	751517	103.69	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	444840	142.84	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	1788973	101.75	ng	0.00
78) Terphenyl-d14	19.60	244	3282114	98.22	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.33	237	18411	9.941	ng	94
53) 2,4-Dinitrophenol	14.39	184	6394	10.190	ng	89
69) Pentachlorophenol	16.63	266	23013	10.202	ng	93

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

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014188.D
Acq On : 08 Feb 2018 11:26
Operator : SJ/JU
Sample : J1161-09
Misc : LOD 8 PPM
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MDL-WATER-03-QT1-2018

Quant Time: Feb 08 17:12:32 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

