

Data Path : Z:\HPCHEM1\BNA N\DATA\BN021518\
 Data File : BN000104.D
 Acq On : 15 Feb 2018 19:17
 Operator : JU/SJ
 Sample : J1161-07
 Misc : 8 PPM LOD
 ALS Vial : 6 Sample Multiplier: 1

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

Quant Time: Feb 16 06:51:24 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 18:24:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	383833	20.00	ng	0.00
21) Naphthalene-d8	11.30	136	1961754	20.00	ng	-0.01
38) Acenaphthene-d10	15.07	164	1184279	20.00	ng	-0.01
63) Phenanthrene-d10	17.79	188	2563896	20.00	ng	-0.01
75) Chrysene-d12	21.90	240	2747187	20.00	ng	-0.02
86) Perylene-d12	24.39	264	2862669	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.93	112	3101976	131.71	ng	-0.01
7) Phenol-d6	7.58	99	4552398	127.61	ng	-0.01
23) Nitrobenzene-d5	9.64	82	3227105	97.81	ng	-0.01
41) 2,4,6-Tribromophenol	16.54	330	1542521	131.59	ng	-0.01
44) 2-Fluorobiphenyl	13.70	172	7892957	93.93	ng	-0.01
78) Terphenyl-d14	20.35	244	9515905	93.53	ng	-0.01
Target Compounds						
40) Hexachlorocyclopentadiene	13.26	237	86728	5.952	ng	92
53) 2,4-Dinitrophenol	15.16	184	25206	12.304	ng	# 14
69) Pentachlorophenol	17.43	266	85186	9.859	ng	94

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

Data Path : Z:\HPCHEM1\BNA N\DATA\BN021518\
Data File : BN000104.D
Acq On : 15 Feb 2018 19:17
Operator : JU/SJ
Sample : J1161-07
Misc : 8 PPM LOD
ALS Vial : 6 Sample Multiplier: 1

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

Quant Time: Feb 16 06:51:24 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 07 18:24:45 2018
Response via : Initial Calibration

