

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080818\
 Data File : BF108030.D
 Acq On : 8 Aug 2018 20:08
 Operator : JU/SJ
 Sample : J3762-07
 Misc : LOD 8 PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-WATER-01-QT3-2018

Quant Time: Aug 09 07:32:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	267892	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1069695	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	575689	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	1193199	20.00	ng	0.00
75) Chrysene-d12	14.21	240	1030745	20.00	ng	-0.01
86) Perylene-d12	15.77	264	900597	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	2052881	138.74	ng	0.01
7) Phenol-d6	6.63	99	2729988	138.84	ng	0.00
23) Nitrobenzene-d5	7.58	82	1498949	78.19	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	898179	156.41	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	2776079	77.42	ng	0.00
78) Terphenyl-d14	13.15	244	3220071	79.43	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	9.16	237	94814	8.303	ng	94
53) 2,4-Dinitrophenol	10.10	184	14070	11.108	ng	# 1
69) Pentachlorophenol	11.35	266	46122	7.063	ng	96

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080818\
Data File : BF108030.D
Acq On : 8 Aug 2018 20:08
Operator : JU/SJ
Sample : J3762-07
Misc : LOD 8 PPM
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LOD-MDL-WATER-01-QT3-2018

Quant Time: Aug 09 07:32:00 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 12:24:24 2018
Response via : Initial Calibration

