

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071317\
 Data File : BF096735.D
 Acq On : 13 Jul 2017 18:44
 Operator : SJ/JU
 Sample : I3757-05
 Misc : 8PPM LOD
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-WATER-05-QT3-2017

Quant Time: Jul 14 17:23:34 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 14:49:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	101553	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	423505	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	192714	20.00	ng	0.00
63) Phenanthrene-d10	11.14	188	326368	20.00	ng	0.00
75) Chrysene-d12	13.77	240	265649	20.00	ng	0.00
86) Perylene-d12	15.15	264	188031	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	827409	122.51	ng	0.00
7) Phenol-d6	6.28	99	1004716	123.96	ng	0.00
23) Nitrobenzene-d5	7.20	82	657893	96.22	ng	0.00
41) 2,4,6-Tribromophenol	10.46	330	213788	129.96	ng	0.00
44) 2-Fluorobiphenyl	9.00	172	1130720	92.70	ng	0.00
78) Terphenyl-d14	12.73	244	1019646	66.57	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	8.79	237	9026	9.303	ng	99
53) 2,4-Dinitrophenol	9.73	184	6123	10.344	ng	# 79
69) Pentachlorophenol	10.95	266	14379	8.020	ng	95

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

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