

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071317\
 Data File : BM010811.D
 Acq On : 13 Jul 2017 14:18
 Operator : SJ/JU
 Sample : I3757-05
 Misc : 8 PPM LOD
 ALS Vial : 11 Sample Multiplier: 1

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

Quant Time: Jul 13 16:15:41 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM071217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 12 14:21:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	169347	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	631934	20.00	ng	0.00
38) Acenaphthene-d10	14.10	164	414708	20.00	ng	0.00
63) Phenanthrene-d10	16.85	188	995630	20.00	ng	0.00
75) Chrysene-d12	21.07	240	1197100	20.00	ng	0.00
86) Perylene-d12	23.20	264	1194799	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.08	112	1392095	136.17	ng	0.00
7) Phenol-d6	6.64	99	1837176	130.36	ng	0.00
23) Nitrobenzene-d5	8.59	82	1679039	102.13	ng	0.00
41) 2,4,6-Tribromophenol	15.60	330	921983	145.48	ng	0.00
44) 2-Fluorobiphenyl	12.71	172	3407094	102.20	ng	0.00
78) Terphenyl-d14	19.51	244	5610107	90.64	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.24	237	78926	7.616	ng	97
53) 2,4-Dinitrophenol	14.22	184	31648	6.616	ng	# 78
69) Pentachlorophenol	16.50	266	71699	7.752	ng	94

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

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