

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071317\
 Data File : BF096734.D
 Acq On : 13 Jul 2017 18:16
 Operator : SJ/JU
 Sample : I3757-05
 Misc : 5PPM LOD
 ALS Vial : 17 Sample Multiplier: 1

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

Quant Time: Jul 14 17:18:09 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	98715	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	413640	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	187221	20.00	ng	0.00
63) Phenanthrene-d10	11.14	188	311435	20.00	ng	0.00
75) Chrysene-d12	13.77	240	208935	20.00	ng	0.00
86) Perylene-d12	15.15	264	150493	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	874247	133.16	ng	0.00
7) Phenol-d6	6.28	99	1053603	133.73	ng	0.00
23) Nitrobenzene-d5	7.20	82	684297	102.47	ng	0.00
41) 2,4,6-Tribromophenol	10.46	330	216787	135.65	ng	0.00
44) 2-Fluorobiphenyl	9.00	172	1169696	98.71	ng	0.00
78) Terphenyl-d14	12.73	244	903428	74.99	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.25	88	19593	5.800	ng	# 84
3) Pyridine	3.02	79	36801	5.184	ng	# 56
4) n-Nitrosodimethylamine	2.85	42	21522	5.421	ng	# 76
32) Benzoic acid	7.65	122	17958	4.447	ng	91
33) 4-Chloroaniline	7.99	127	42272	5.445	ng	96
35) Caprolactam	8.32	113	9178	5.009	ng	# 57
52) 3-Nitroaniline	9.61	138	18550	5.128	ng	89
55) 4-Nitrophenol	9.78	139	11321	4.284	ng	# 60
64) 4,6-Dinitro-2-methylphenol	10.26	198	6958	3.558	ng	# 77

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

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