

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
 Data File : BM016043.D
 Acq On : 24 Jul 2018 01:45
 Operator : SJ/JU
 Sample : J3762-07
 Misc : LOD 5 PPM
 ALS Vial : 16 Sample Multiplier: 1

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

Quant Time: Jul 24 08:26:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 23 16:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	38932	20.00	ng	0.00
21) Naphthalene-d8	11.05	136	148518	20.00	ng	0.00
38) Acenaphthene-d10	14.85	164	78745	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	171272	20.00	ng	0.00
75) Chrysene-d12	21.70	240	188564	20.00	ng	0.00
86) Perylene-d12	24.07	264	201928	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	301162	128.50	ng	0.00
7) Phenol-d6	7.39	99	372661	121.75	ng	0.00
23) Nitrobenzene-d5	9.42	82	284013	88.94	ng	0.00
41) 2,4,6-Tribromophenol	16.34	330	129901	130.06	ng	0.00
44) 2-Fluorobiphenyl	13.48	172	570386	88.13	ng	0.00
78) Terphenyl-d14	20.15	244	1087888	120.77	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.52	88	4348	3.960	ng	# 100
3) Pyridine	4.01	79	6960	2.631	ng	# 61
4) n-Nitrosodimethylamine	3.90	42	7911	3.834	ng	# 34
32) Benzoic acid	10.35	122	3310	2.323	ng	# 92
33) 4-Chloroaniline	11.25	127	7359	2.399	ng	# 87
35) Caprolactam	12.01	113	1142	1.856	ng	# 68
52) 3-Nitroaniline	14.80	138	3459	2.310	ng	# 72
55) 4-Nitrophenol	15.13	139	1740	1.621	ng	# 79
64) 4,6-Dinitro-2-methylphenol	16.00	198	2705	2.602	ng	# 86

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

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