

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
 Data File : BM016051.D
 Acq On : 24 Jul 2018 06:39
 Operator : SJ/JU
 Sample : J3762-09
 Misc : MDL 5 PPM
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-03-QT3-2018

Quant Time: Jul 24 09:36:21 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	39581	20.00	ng	0.00
21) Naphthalene-d8	11.05	136	148661	20.00	ng	0.00
38) Acenaphthene-d10	14.85	164	81499	20.00	ng	0.00
63) Phenanthrene-d10	17.58	188	171644	20.00	ng	0.00
75) Chrysene-d12	21.70	240	191629	20.00	ng	0.00
86) Perylene-d12	24.07	264	205621	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.74	112	297860	125.00	ng	0.00
7) Phenol-d6	7.38	99	371279	119.31	ng	0.00
23) Nitrobenzene-d5	9.41	82	286629	89.68	ng	0.00
41) 2,4,6-Tribromophenol	16.33	330	130907	126.64	ng	0.00
44) 2-Fluorobiphenyl	13.47	172	580795	86.71	ng	0.00
78) Terphenyl-d14	20.15	244	1108853	121.13	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	4706	4.215	ng	# 100
3) Pyridine	4.00	79	7090	2.636	ng	# 79
4) n-Nitrosodimethylamine	3.90	42	6478	3.088	ng	# 38
32) Benzoic acid	10.43	122	115	0.081	ng	# 85
33) 4-Chloroaniline	11.25	127	7722	2.515	ng	# 86
35) Caprolactam	12.00	113	1430	2.322	ng	# 62
52) 3-Nitroaniline	14.80	138	3661	2.362	ng	# 85
55) 4-Nitrophenol	15.12	139	1849	1.664	ng	# 62
64) 4,6-Dinitro-2-methylphenol	16.00	198	2423	2.326	ng	# 77

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

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