

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
 Data File : BM014187.D
 Acq On : 08 Feb 2018 10:50
 Operator : SJ/JU
 Sample : J1161-09
 Misc : LOD 5 PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-03-QT1-2018

Quant Time: Feb 08 17:07:30 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 17:16:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	64392	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	294387	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	204661	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	533514	20.00	ng	0.00
75) Chrysene-d12	21.16	240	634391	20.00	ng	0.00
86) Perylene-d12	23.34	264	606436	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	528311	143.00	ng	0.00
7) Phenol-d6	6.73	99	716262	136.89	ng	0.00
23) Nitrobenzene-d5	8.70	82	634393	96.26	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	398808	137.33	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	1475499	90.00	ng	0.00
78) Terphenyl-d14	19.60	244	2799185	86.22	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.15	88	7797	5.136	ng	# 100
3) Pyridine	3.55	79	16163	3.845	ng	# 89
4) n-Nitrosodimethylamine	3.46	42	13319	4.974	ng	# 41
32) Benzoic acid	9.75	122	1075	6.464	ng	# 66
33) 4-Chloroaniline	10.52	127	22542	3.364	ng	# 84
35) Caprolactam	11.30	113	4307	2.763	ng	# 48
52) 3-Nitroaniline	14.15	138	13737	3.532	ng	# 78
55) 4-Nitrophenol	14.52	139	3879	6.960	ng	# 61
64) 4,6-Dinitro-2-methylphenol	15.39	198	8555	7.638	ng	# 74

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

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