

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020718\
 Data File : BM014178.D
 Acq On : 08 Feb 2018 01:29
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD 5 PPM
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT1-2018

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:50:43 PM

Quant Time: Feb 08 12:18:24 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	7.54	152	79435	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	360631	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	244909	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	631746	20.00	ng	0.00
75) Chrysene-d12	21.16	240	747953	20.00	ng	0.00
86) Perylene-d12	23.34	264	694943	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	625875	137.33	ng	0.00
7) Phenol-d6	6.73	99	855549	132.55	ng	0.00
23) Nitrobenzene-d5	8.70	82	769467	95.31	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	496816	142.96	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	1795603	91.52	ng	0.00
78) Terphenyl-d14	19.60	244	3339969	87.25	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.15	88	9079	4.85	ng	Qvalue # 100
3) Pyridine	3.54	79	19274	3.72	ng	90
4) n-Nitrosodimethylamine	3.46	42	14168	4.29	ng	# 51
32) Benzoic acid	9.65	122	8773m	8.04	ng	
33) 4-Chloroaniline	10.52	127	26816	3.27	ng	# 86
35) Caprolactam	11.30	113	5768m	3.02	ng	
52) 3-Nitroaniline	14.15	138	15290	3.29	ng	# 58
55) 4-Nitrophenol	14.52	139	6551m	7.44	ng	
64) 4,6-Dinitro-2-methylphenol	15.39	198	10315	7.68	ng	80

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

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