

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102418\
 Data File : BN003286.D
 Acq On : 24 Oct 2018 23:49
 Operator : MJ/SJ
 Sample : J5164-07
 Misc : LOD 5PPM
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2018

Quant Time: Oct 25 07:08:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 24 14:21:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	93439	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	426170	20.00	ng	0.00
39) Acenaphthene-d10	14.24	164	272083	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	638934	20.00	ng	0.00
76) Chrysene-d12	21.21	240	725289	20.00	ng	0.00
87) Perylene-d12	23.41	264	757836	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	728815	135.85	ng	0.00
7) Phenol-d6	6.80	99	1010580	132.48	ng	0.00
23) Nitrobenzene-d5	8.76	82	596611	80.29	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	497300	135.74	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1595174	79.52	ng	0.00
79) Terphenyl-d14	19.64	244	2575716	75.95	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.21	88	10628	4.172	ng	Qvalue 85
3) Pyridine	3.61	79	24100	4.140	ng	93
4) n-Nitrosodimethylamine	3.53	42	8415	3.927	ng	# 96
32) Benzoic acid	9.75	122	3303	5.622	ng	# 80
33) 4-Chloroaniline	10.57	127	36316	3.946	ng	98
35) Caprolactam	11.33	113	8320	3.318	ng	93
53) 3-Nitroaniline	14.20	138	18393	3.446	ng	# 94
56) 4-Nitrophenol	14.65	139	55429	18.851	ng	# 18
65) 4,6-Dinitro-2-methylphenol	15.45	198	12124	2.719	ng	99

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

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