

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022318\
 Data File : BF103196.D
 Acq On : 23 Feb 2018 13:45
 Operator : SJ/JU
 Sample : J1161-07
 Misc : LOD-S-5ppm
 ALS Vial : 14 Sample Multiplier: 1

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

Quant Time: Feb 23 17:02:35 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	175913	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	768428	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	354695	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	610832	20.00	ng	0.00
75) Chrysene-d12	14.04	240	453083	20.00	ng	0.00
86) Perylene-d12	15.54	264	438017	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1318215	113.33	ng	0.00
7) Phenol-d6	6.50	99	1596300	112.16	ng	0.00
23) Nitrobenzene-d5	7.44	82	1207957	89.77	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	356488	118.74	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1771526	86.35	ng	0.00
78) Terphenyl-d14	12.99	244	1453266	77.10	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.65	88	29011	4.722	ng	96
3) Pyridine	3.56	79	59632	3.636	ng	96
4) n-Nitrosodimethylamine	3.40	42	18215	2.754	ng	# 95
32) Benzoic acid	7.88	122	22462	9.345	ng	90
33) 4-Chloroaniline	8.23	127	79472	4.895	ng	97
35) Caprolactam	8.56	113	16906	4.713	ng	# 72
52) 3-Nitroaniline	9.86	138	35350	4.666	ng	93
55) 4-Nitrophenol	10.04	139	15592	3.321	ng	# 82
64) 4,6-Dinitro-2-methylphenol	10.51	198	11411	7.699	ng	84

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

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