

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102318\
 Data File : BF110093.D
 Acq On : 23 Oct 2018 22:27
 Operator : JU/SJ
 Sample : J5164-07
 Misc : LOD 5PPM
 ALS Vial : 15 Sample Multiplier: 1

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

Quant Time: Oct 24 07:30:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	163201	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	693261	20.00	ng	-0.02
39) Acenaphthene-d10	10.00	164	341455	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	653306	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	560446	20.00	ng	-0.02
87) Perylene-d12	15.67	264	484764	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	1235913	123.32	ng	-0.02
7) Phenol-d6	6.59	99	1543712	120.43	ng	-0.02
23) Nitrobenzene-d5	7.53	82	935870	80.07	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	412711	122.02	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	1724009	79.28	ng	-0.02
79) Terphenyl-d14	13.09	244	1820479	67.55	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	2.83	88	21837	4.159	ng	91
3) Pyridine	3.66	79	54713	4.678	ng	# 82
4) n-Nitrosodimethylamine	3.55	42	24864	5.075	ng	# 88
32) Benzoic acid	7.94	122	16004	2.175	ng	88
33) 4-Chloroaniline	8.32	127	55543	3.863	ng	96
35) Caprolactam	8.65	113	15381	4.588	ng	# 86
53) 3-Nitroaniline	9.95	138	26949	4.174	ng	92
56) 4-Nitrophenol	10.10	139	22582	4.726	ng	92
65) 4,6-Dinitro-2-methylphenol	10.59	198	9454	3.167	ng	88

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

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