

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
 Data File : BF107507.D
 Acq On : 19 Jul 2018 10:41
 Operator : JU/SJ
 Sample : J4006-01DL 20X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RR-BASELINE-WATERDL

Quant Time: Jul 19 13:26:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 11:55:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	99817	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	371185	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	201664	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	375135	20.00	ng	0.00
75) Chrysene-d12	14.17	240	282923	20.00	ng	-0.01
86) Perylene-d12	15.70	264	241069	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	17454	2.66	ng	-0.01
7) Phenol-d6	6.60	99	13790	1.62	ng	-0.01
23) Nitrobenzene-d5	7.54	82	33824	4.86	ng	-0.02
41) 2,4,6-Tribromophenol	10.81	330	12807	7.23	ng	-0.02
44) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
78) Terphenyl-d14	13.10	244	66247	5.82	ng	-0.01
Target Compounds						
17) 2-Methylphenol	7.22	107	16132	2.553	ng	# 86
20) 3+4-Methylphenols	7.38	107	77463	9.650	ng	# 90
27) 2,4-Dimethylphenol	7.92	122	106776	19.256	ng	# 78

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

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