

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061118\
 Data File : BF106324.D
 Acq On : 12 Jun 2018 5:31
 Operator : JU/SJ
 Sample : J3311-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-35

Quant Time: Jun 12 09:44:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	197484	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	690377	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	309603	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	513808	20.00	ng	0.00
75) Chrysene-d12	14.14	240	502487	20.00	ng	-0.02
86) Perylene-d12	15.67	264	423675	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	731229	62.67	ng	0.00
7) Phenol-d6	6.58	99	606656	41.15	ng	0.00
23) Nitrobenzene-d5	7.54	82	1098843	93.63	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	454653	152.62	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1765673	111.32	ng	0.00
78) Terphenyl-d14	13.09	244	1595865	78.88	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	7.36	107	31448	2.197	ng	87
49) Dimethylphthalate	9.72	163	86565	3.914	ng	# 90
51) Acenaphthene	10.05	154	43053	2.270	ng	96

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

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