

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121823\
 Data File : BG060073.D
 Acq On : 18 Dec 2023 19:35
 Operator : MA/JU
 Sample : 05592-01DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW04SDL

Quant Time: Dec 19 04:41:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 19 04:37:23 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	65021	20.000	ng	0.00
21) Naphthalene-d8	10.755	136	258669	20.000	ng	# 0.00
39) Acenaphthene-d10	14.591	164	216064	20.000	ng	0.00
64) Phenanthrene-d10	17.341	188	642189	20.000	ng	# 0.00
76) Chrysene-d12	21.613	240	709895	20.000	ng	# 0.00
86) Perylene-d12	24.803	264	811284	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.520	112	36864	9.705	ng	0.00
7) Phenol-d6	7.106	99	31782	5.948	ng	0.00
23) Nitrobenzene-d5	9.110	82	87245	15.315	ng	-0.01
42) 2,4,6-Tribromophenol	16.078	330	129418	25.702	ng	0.00
45) 2-Fluorobiphenyl	13.211	172	300409	17.017	ng	0.00
79) Terphenyl-d14	19.962	244	775055	19.933	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	12.635	142	82189	7.878	ng	97
52) Acenaphthene	14.656	154	314132	24.041	ng	95
58) Fluorene	15.637	166	98673	5.502	ng	96

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

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