

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071222\
 Data File : BG054258.D
 Acq On : 12 Jul 2022 16:11
 Operator : CG/JU
 Sample : PB146157BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB146157BL

Quant Time: Jul 13 01:36:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 15:49:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.038	152	18395	20.000	ng	# 0.00
21) Naphthalene-d8	10.846	136	73659	20.000	ng	# 0.00
39) Acenaphthene-d10	14.671	164	59310	20.000	ng	0.00
64) Phenanthrene-d10	17.415	188	173773	20.000	ng	# 0.00
76) Chrysene-d12	21.686	240	236151	20.000	ng	# 0.00
86) Perylene-d12	24.924	264	234753	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.611	112	104891	114.562	ng	0.00
7) Phenol-d6	7.209	99	141009	111.662	ng	0.00
23) Nitrobenzene-d5	9.201	82	96018	70.739	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	148726	129.452	ng	0.00
45) 2-Fluorobiphenyl	13.296	172	318600	78.930	ng	0.00
79) Terphenyl-d14	20.024	244	996162	86.745	ng	0.00

Target Compounds	Qvalue

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

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