

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082022\
 Data File : BG054680.D
 Acq On : 20 Aug 2022 3:31
 Operator : CG/JU
 Sample : PB147107BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB147107BL

Quant Time: Aug 22 00:38:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 00:08:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.161	152	57574	20.000	ng	0.00
21) Naphthalene-d8	10.981	136	231717	20.000	ng	0.00
39) Acenaphthene-d10	14.789	164	150273	20.000	ng	0.00
64) Phenanthrene-d10	17.532	188	342290	20.000	ng	0.00
76) Chrysene-d12	21.827	240	339921	20.000	ng	0.00
86) Perylene-d12	25.194	264	343258	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.693	112	438339	137.859	ng	0.00
7) Phenol-d6	7.297	99	577024	129.763	ng	0.00
23) Nitrobenzene-d5	9.330	82	341605	85.113	ng	0.00
42) 2,4,6-Tribromophenol	16.275	330	197683	125.986	ng	0.00
45) 2-Fluorobiphenyl	13.414	172	813976	84.846	ng	0.00
79) Terphenyl-d14	20.135	244	1364708	85.810	ng	0.00

Target Compounds	Qvalue

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

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