

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092222\
 Data File : BG055008.D
 Acq On : 22 Sep 2022 13:19
 Operator : CG/JU
 Sample : PB147786BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB147786BL

Quant Time: Sep 23 03:26:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 17:41:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	40732	20.000	ng	0.00
21) Naphthalene-d8	10.885	136	163253	20.000	ng	# 0.00
39) Acenaphthene-d10	14.704	164	111785	20.000	ng	0.00
64) Phenanthrene-d10	17.448	188	277413	20.000	ng	0.00
76) Chrysene-d12	21.725	240	308886	20.000	ng	0.00
86) Perylene-d12	25.033	264	323624	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.591	112	307640	147.784	ng	0.00
7) Phenol-d6	7.219	99	385498	138.160	ng	0.00
23) Nitrobenzene-d5	9.234	82	223295	79.279	ng	0.00
42) 2,4,6-Tribromophenol	16.191	330	172664	131.849	ng	0.00
45) 2-Fluorobiphenyl	13.323	172	597369	84.219	ng	0.00
79) Terphenyl-d14	20.051	244	1218597	82.525	ng	0.00

Target Compounds Qvalue

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

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