

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092322\
 Data File : BG055014.D
 Acq On : 23 Sep 2022 16:02
 Operator : CG/JU
 Sample : N4751-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-7

Quant Time: Sep 24 02:28:00 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	8.056	152	38974	20.000	ng	0.00
21) Naphthalene-d8	10.882	136	153527	20.000	ng	# 0.00
39) Acenaphthene-d10	14.701	164	108559	20.000	ng	0.00
64) Phenanthrene-d10	17.451	188	265701	20.000	ng	# 0.00
76) Chrysene-d12	21.728	240	270854	20.000	ng	0.00
86) Perylene-d12	25.036	264	285863	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.600	112	226695	113.812	ng	0.00
7) Phenol-d6	7.221	99	295398	110.644	ng	0.00
23) Nitrobenzene-d5	9.237	82	168315	63.545	ng	0.00
42) 2,4,6-Tribromophenol	16.193	330	134994	106.147	ng	0.00
45) 2-Fluorobiphenyl	13.326	172	468810	68.058	ng	0.00
79) Terphenyl-d14	20.053	244	913284	70.533	ng	0.00

Target Compounds	Qvalue

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

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