

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071222\
 Data File : BG054266.D
 Acq On : 12 Jul 2022 21:54
 Operator : CG/JU
 Sample : N3649-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B-2-UC

Quant Time: Jul 13 01:37:07 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	8.036	152	14175	20.000	ng	# 0.00
21) Naphthalene-d8	10.851	136	54768	20.000	ng	# 0.00
39) Acenaphthene-d10	14.670	164	46454	20.000	ng	0.00
64) Phenanthrene-d10	17.414	188	144979	20.000	ng	# 0.00
76) Chrysene-d12	21.685	240	183430	20.000	ng	# 0.00
86) Perylene-d12	24.917	264	187518	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.610	112	86611	122.758	ng	0.00
7) Phenol-d6	7.208	99	113115	116.240	ng	0.00
23) Nitrobenzene-d5	9.200	82	74639	73.956	ng	0.00
42) 2,4,6-Tribromophenol	16.156	330	128374	142.660	ng	0.00
45) 2-Fluorobiphenyl	13.289	172	244507	77.338	ng	0.00
79) Terphenyl-d14	20.022	244	664769	74.525	ng	0.00

Target Compounds Qvalue

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

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