

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080522\
 Data File : BG054531.D
 Acq On : 5 Aug 2022 11:50
 Operator : CG/JU
 Sample : PB146695BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB146695BL

Quant Time: Aug 06 05:39:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.973	152	15486	20.000	ng	#-0.02
21) Naphthalene-d8	10.787	136	59856	20.000	ng	# 0.00
39) Acenaphthene-d10	14.618	164	47334	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	132819	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	151505	20.000	ng	# 0.00
86) Perylene-d12	24.841	264	157802	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.564	112	110413	148.863	ng	0.00
7) Phenol-d6	7.180	99	144649	142.326	ng	0.00
23) Nitrobenzene-d5	9.142	82	95023	86.452	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	121501	122.254	ng	0.00
45) 2-Fluorobiphenyl	13.237	172	288070	85.853	ng	0.00
79) Terphenyl-d14	19.976	244	720515	94.344	ng	0.00

Target Compounds Qvalue

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

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