

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021122\
 Data File : BG052367.D
 Acq On : 11 Feb 2022 11:15
 Operator : CG/JU
 Sample : PB142428BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB142428BL

Quant Time: Feb 11 13:50:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	23058	20.000	ng	0.00
21) Naphthalene-d8	11.067	136	96025	20.000	ng	# 0.00
39) Acenaphthene-d10	14.873	164	70096	20.000	ng	0.00
64) Phenanthrene-d10	17.622	188	188009	20.000	ng	0.00
76) Chrysene-d12	21.940	240	222305	20.000	ng	0.00
86) Perylene-d12	25.412	264	212106	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.762	112	188531	142.503	ng	0.01
7) Phenol-d6	7.378	99	261244	127.979	ng	0.00
23) Nitrobenzene-d5	9.404	82	180097	81.512	ng	0.00
42) 2,4,6-Tribromophenol	16.365	330	143807	142.802	ng	0.00
45) 2-Fluorobiphenyl	13.499	172	444583	88.131	ng	0.00
79) Terphenyl-d14	20.219	244	1040495	82.651	ng	0.00

Target Compounds	Qvalue

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

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