

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042322\
 Data File : BG053287.D
 Acq On : 23 Apr 2022 11:12
 Operator : CG/JU
 Sample : PB144287BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB144287BL

Quant Time: Apr 24 02:48:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.108	152	29050	20.000	ng	0.00
21) Naphthalene-d8	10.921	136	114700	20.000	ng	# 0.00
39) Acenaphthene-d10	14.728	164	83755	20.000	ng	0.00
64) Phenanthrene-d10	17.477	188	213433	20.000	ng	0.00
76) Chrysene-d12	21.760	240	238588	20.000	ng	0.00
86) Perylene-d12	25.049	264	231361	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.676	112	227841	134.444	ng	0.00
7) Phenol-d6	7.268	99	319667	122.319	ng	0.00
23) Nitrobenzene-d5	9.271	82	229789	81.345	ng	0.00
42) 2,4,6-Tribromophenol	16.220	330	159717	119.862	ng	0.00
45) 2-Fluorobiphenyl	13.353	172	512042	84.300	ng	0.00
79) Terphenyl-d14	20.080	244	1119069	79.216	ng	0.00

Target Compounds	Qvalue

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

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