

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042922\
 Data File : BG053376.D
 Acq On : 29 Apr 2022 14:17
 Operator : CG/JU
 Sample : PB144441BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB144441BL

Quant Time: May 02 01:02:38 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	30222	20.000	ng	0.00
21) Naphthalene-d8	10.910	136	116087	20.000	ng	-0.01
39) Acenaphthene-d10	14.722	164	84384	20.000	ng	-0.01
64) Phenanthrene-d10	17.472	188	211200	20.000	ng	-0.01
76) Chrysene-d12	21.754	240	246355	20.000	ng	-0.01
86) Perylene-d12	25.050	264	238856	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.670	112	244107	138.457	ng	0.00
7) Phenol-d6	7.262	99	336501	123.768	ng	-0.01
23) Nitrobenzene-d5	9.265	82	239858	83.894	ng	-0.01
42) 2,4,6-Tribromophenol	16.215	330	161090	119.991	ng	-0.01
45) 2-Fluorobiphenyl	13.348	172	531801	86.901	ng	-0.01
79) Terphenyl-d14	20.074	244	1157491	79.353	ng	-0.01

Target Compounds	Qvalue

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

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