

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061022\
 Data File : BG053873.D
 Acq On : 10 Jun 2022 13:36
 Operator : CG/JU
 Sample : PB145428BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB145428BL

Quant Time: Jun 10 17:33:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 02 23:22:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.090	152	23379	20.000	ng	0.00
21) Naphthalene-d8	10.898	136	90854	20.000	ng	# 0.00
39) Acenaphthene-d10	14.711	164	65516	20.000	ng	0.00
64) Phenanthrene-d10	17.455	188	172706	20.000	ng	# 0.00
76) Chrysene-d12	21.732	240	208929	20.000	ng	0.00
86) Perylene-d12	25.005	264	209031	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.657	112	173978	136.655	ng	0.00
7) Phenol-d6	7.244	99	224141	129.969	ng	0.00
23) Nitrobenzene-d5	9.247	82	144229	104.247	ng	0.00
42) 2,4,6-Tribromophenol	16.198	330	131730	165.641	ng	0.00
45) 2-Fluorobiphenyl	13.342	172	393635	89.884	ng	0.00
79) Terphenyl-d14	20.064	244	926497	88.220	ng	0.00

Target Compounds	Qvalue

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

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