

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102122\
 Data File : BG055238.D
 Acq On : 21 Oct 2022 17:44
 Operator : CG/JU
 Sample : N5217-18
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-C1-COMP

Quant Time: Oct 22 01:30:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 09:21:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.284	152	39197	20.000	ng	0.00
21) Naphthalene-d8	11.116	136	179982	20.000	ng	0.00
39) Acenaphthene-d10	14.894	164	129749	20.000	ng	0.00
64) Phenanthrene-d10	17.626	188	306467	20.000	ng	0.00
76) Chrysene-d12	21.910	240	235979	20.000	ng	0.00
86) Perylene-d12	25.317	264	217183	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.811	112	331597	152.203	ng	0.00
7) Phenol-d6	7.427	99	465075	140.092	ng	0.00
23) Nitrobenzene-d5	9.459	82	290209	77.845	ng	0.00
42) 2,4,6-Tribromophenol	16.375	330	196488	146.146	ng	0.00
45) 2-Fluorobiphenyl	13.525	172	653223	80.417	ng	0.00
79) Terphenyl-d14	20.206	244	1085592	85.719	ng	0.00

Target Compounds Qvalue

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

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