

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102622\
 Data File : BG055311.D
 Acq On : 26 Oct 2022 23:25
 Operator : CG/JU
 Sample : N5218-31
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-4

Quant Time: Oct 27 10:44:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 24 16:54:05 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.267	152	33793	20.000	ng	0.00
21) Naphthalene-d8	11.099	136	145794	20.000	ng	0.00
39) Acenaphthene-d10	14.883	164	99058	20.000	ng	0.00
64) Phenanthrene-d10	17.615	188	222743	20.000	ng	0.00
76) Chrysene-d12	21.893	240	212796	20.000	ng	0.00
86) Perylene-d12	25.289	264	227498	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.794	112	222764	115.424	ng	0.00
7) Phenol-d6	7.415	99	325831	116.532	ng	0.00
23) Nitrobenzene-d5	9.442	82	194774	68.218	ng	0.00
42) 2,4,6-Tribromophenol	16.364	330	128159	119.013	ng	0.00
45) 2-Fluorobiphenyl	13.508	172	442117	73.052	ng	0.00
79) Terphenyl-d14	20.189	244	724918	70.779	ng	0.00

Target Compounds Qvalue

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

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