

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
 Data File : BG055593.D
 Acq On : 12 Nov 2022 17:19
 Operator : CG/JU
 Sample : N5431-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221178-COMP-13-MODIFIED-ELUTRIATE

Quant Time: Nov 14 03:28:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 03:39:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.267	152	60953	20.000	ng	0.00
21) Naphthalene-d8	11.093	136	255007	20.000	ng	0.00
39) Acenaphthene-d10	14.888	164	190903	20.000	ng	0.00
64) Phenanthrene-d10	17.626	188	478982	20.000	ng	0.00
76) Chrysene-d12	21.927	240	470517	20.000	ng	0.00
86) Perylene-d12	25.347	264	493707	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.793	112	157009	49.064	ng	0.00
7) Phenol-d6	7.397	99	151384	34.247	ng	0.00
23) Nitrobenzene-d5	9.436	82	274957	73.095	ng	0.00
42) 2,4,6-Tribromophenol	16.369	330	246954	132.582	ng	0.00
45) 2-Fluorobiphenyl	13.514	172	680515	53.770	ng	0.00
79) Terphenyl-d14	20.218	244	1269694	54.201	ng	0.00

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055593.D
Acq On : 12 Nov 2022 17:19
Operator : CG/JU
Sample : N5431-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
20221178-COMP-13-MODIFIED-ELUTRIATE

Quant Time: Nov 14 03:28:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 10 03:39:46 2022
Response via : Initial Calibration

