

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110922\
 Data File : BG055531.D
 Acq On : 9 Nov 2022 20:11
 Operator : CG/JU
 Sample : N5360-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221175-COMP-10-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:31:51 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	8.272	152	58398	20.000	ng	0.00
21) Naphthalene-d8	11.098	136	240214	20.000	ng	0.00
39) Acenaphthene-d10	14.888	164	163766	20.000	ng	0.00
64) Phenanthrene-d10	17.626	188	376819	20.000	ng	0.00
76) Chrysene-d12	21.921	240	363355	20.000	ng	0.00
86) Perylene-d12	25.341	264	377182	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.793	112	121647	39.677	ng	0.00
7) Phenol-d6	7.397	99	147930	34.930	ng	0.00
23) Nitrobenzene-d5	9.436	82	273230	77.109	ng	0.00
42) 2,4,6-Tribromophenol	16.363	330	126316	79.053	ng	0.00
45) 2-Fluorobiphenyl	13.519	172	637971	58.761	ng	0.00
79) Terphenyl-d14	20.211	244	454021	25.097	ng	0.00

Target Compounds	Qvalue

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

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