

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
 Data File : BG055591.D
 Acq On : 12 Nov 2022 15:53
 Operator : CG/JU
 Sample : N5431-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221176-COMP-11-MODIFIED-ELUTRIATE

Quant Time: Nov 14 03:27:45 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	8.270	152	79208	20.000	ng	0.00
21) Naphthalene-d8	11.097	136	326431	20.000	ng	0.00
39) Acenaphthene-d10	14.886	164	245451	20.000	ng	0.00
64) Phenanthrene-d10	17.630	188	602941	20.000	ng	0.00
76) Chrysene-d12	21.925	240	603038	20.000	ng	0.00
86) Perylene-d12	25.356	264	645612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	122685	29.502	ng	0.00
7) Phenol-d6	7.401	99	125302	21.813	ng	0.00
23) Nitrobenzene-d5	9.440	82	200179	41.572	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	188598	78.751	ng	0.00
45) 2-Fluorobiphenyl	13.511	172	506295	31.114	ng	0.00
79) Terphenyl-d14	20.215	244	981600	32.694	ng	0.00

Target Compounds	Qvalue

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

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