

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111522\
 Data File : BG055634.D
 Acq On : 15 Nov 2022 23:13
 Operator : CG/JU
 Sample : N5431-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221170-COMP-05-MODIFIED-ELUTRIATE

Quant Time: Nov 16 01:47:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 01:36:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.258	152	40781	20.000	ng	0.00
21) Naphthalene-d8	11.084	136	171682	20.000	ng	0.00
39) Acenaphthene-d10	14.880	164	127624	20.000	ng	0.00
64) Phenanthrene-d10	17.618	188	329198	20.000	ng	0.00
76) Chrysene-d12	21.913	240	314401	20.000	ng	0.00
86) Perylene-d12	25.326	264	329122	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.779	112	92785	43.336	ng	0.00
7) Phenol-d6	7.389	99	100900	34.117	ng	0.00
23) Nitrobenzene-d5	9.427	82	145586	57.487	ng	0.00
42) 2,4,6-Tribromophenol	16.360	330	138268	111.038	ng	0.00
45) 2-Fluorobiphenyl	13.505	172	367777	43.468	ng	0.00
79) Terphenyl-d14	20.203	244	732633	46.804	ng	0.00

Target Compounds	Qvalue

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

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