

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110422\
 Data File : BG055455.D
 Acq On : 4 Nov 2022 15:39
 Operator : CG/JU
 Sample : N5306-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221166-COMP-01-MODIFIED-ELUTRIATE

Quant Time: Nov 05 00:03:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 05:28:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	25656	20.000	ng	0.00
21) Naphthalene-d8	11.061	136	101580	20.000	ng	0.00
39) Acenaphthene-d10	14.851	164	73286	20.000	ng	0.00
64) Phenanthrene-d10	17.583	188	182875	20.000	ng	0.00
76) Chrysene-d12	21.855	240	196941	20.000	ng	0.00
86) Perylene-d12	25.227	264	214969	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.762	112	106106	74.271	ng	0.00
7) Phenol-d6	7.383	99	106580	53.797	ng	0.00
23) Nitrobenzene-d5	9.410	82	155096	81.373	ng	0.00
42) 2,4,6-Tribromophenol	16.332	330	141967	142.018	ng	0.00
45) 2-Fluorobiphenyl	13.476	172	402913	81.514	ng	0.00
79) Terphenyl-d14	20.157	244	796386	78.633	ng	0.00

Target Compounds Qvalue

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

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