

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110422\
 Data File : BG055456.D
 Acq On : 4 Nov 2022 16:20
 Operator : CG/JU
 Sample : N5306-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221174-COMP-09-MODIFIED-ELUTRIATE

Quant Time: Nov 05 00:03:12 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	8.235	152	27284	20.000	ng	0.00
21) Naphthalene-d8	11.061	136	111734	20.000	ng	0.00
39) Acenaphthene-d10	14.851	164	79787	20.000	ng	0.00
64) Phenanthrene-d10	17.583	188	199155	20.000	ng	0.00
76) Chrysene-d12	21.854	240	210316	20.000	ng	0.00
86) Perylene-d12	25.227	264	225308	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.756	112	103032	67.816	ng	0.00
7) Phenol-d6	7.383	99	101786	48.312	ng	0.00
23) Nitrobenzene-d5	9.410	82	158623	75.660	ng	0.00
42) 2,4,6-Tribromophenol	16.331	330	147801	135.807	ng	0.00
45) 2-Fluorobiphenyl	13.476	172	417145	77.517	ng	0.00
79) Terphenyl-d14	20.156	244	763407	70.583	ng	0.00

Target Compounds	Qvalue

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

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