

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110422\
 Data File : BG055454.D
 Acq On : 4 Nov 2022 14:58
 Operator : CG/JU
 Sample : N5306-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221173-COMP-08-MODIFIED-ELUTRIATE

Quant Time: Nov 05 00:02:47 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	25279	20.000	ng	0.00
21) Naphthalene-d8	11.061	136	103475	20.000	ng	0.00
39) Acenaphthene-d10	14.850	164	73686	20.000	ng	0.00
64) Phenanthrene-d10	17.582	188	185989	20.000	ng	0.00
76) Chrysene-d12	21.854	240	200613	20.000	ng	0.00
86) Perylene-d12	25.226	264	215175	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.755	112	110066	78.192	ng	0.00
7) Phenol-d6	7.383	99	105966	54.285	ng	0.00
23) Nitrobenzene-d5	9.410	82	166946	85.986	ng	0.00
42) 2,4,6-Tribromophenol	16.331	330	148988	148.233	ng	0.00
45) 2-Fluorobiphenyl	13.476	172	431010	86.725	ng	0.00
79) Terphenyl-d14	20.156	244	862051	83.558	ng	0.00

Target Compounds	Qvalue

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

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