

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110322\
 Data File : BG055447.D
 Acq On : 3 Nov 2022 21:32
 Operator : CG/JU
 Sample : N5306-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

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

Quant Time: Nov 04 05:31:48 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	8.234	152	31068	20.000	ng	0.00
21) Naphthalene-d8	11.066	136	125408	20.000	ng	0.00
39) Acenaphthene-d10	14.850	164	88365	20.000	ng	0.00
64) Phenanthrene-d10	17.582	188	207529	20.000	ng	0.00
76) Chrysene-d12	21.854	240	207308	20.000	ng	0.00
86) Perylene-d12	25.226	264	223851	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.761	112	114099	65.954	ng	0.00
7) Phenol-d6	7.382	99	114662	47.795	ng	0.00
23) Nitrobenzene-d5	9.409	82	193085	82.056	ng	0.00
42) 2,4,6-Tribromophenol	16.331	330	150025	124.469	ng	0.00
45) 2-Fluorobiphenyl	13.481	172	488306	81.932	ng	0.00
79) Terphenyl-d14	20.161	244	839344	78.730	ng	0.00

Target Compounds	Qvalue

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

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