

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
 Data File : BG055594.D
 Acq On : 12 Nov 2022 18:01
 Operator : CG/JU
 Sample : N5431-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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

Quant Time: Nov 14 03:29:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 03:39:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.266	152	55689	20.000	ng	0.00
21) Naphthalene-d8	11.098	136	234240	20.000	ng	0.00
39) Acenaphthene-d10	14.888	164	176329	20.000	ng	0.00
64) Phenanthrene-d10	17.626	188	445500	20.000	ng	0.00
76) Chrysene-d12	21.927	240	452350	20.000	ng	0.00
86) Perylene-d12	25.346	264	483292	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.793	112	84420	28.874	ng	0.00
7) Phenol-d6	7.397	99	87190	21.589	ng	0.00
23) Nitrobenzene-d5	9.436	82	147366	42.649	ng	0.00
42) 2,4,6-Tribromophenol	16.363	330	140993	81.951	ng	0.00
45) 2-Fluorobiphenyl	13.513	172	378951	32.417	ng	0.00
79) Terphenyl-d14	20.211	244	727848	32.318	ng	0.00

Target Compounds	Qvalue

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

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