

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110922\
 Data File : BG055529.D
 Acq On : 9 Nov 2022 18:48
 Operator : CG/JU
 Sample : N5360-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221169-COMP-04-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:31:01 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	8.271	152	56155	20.000	ng	0.00
21) Naphthalene-d8	11.097	136	227120	20.000	ng	0.00
39) Acenaphthene-d10	14.887	164	152989	20.000	ng	0.00
64) Phenanthrene-d10	17.625	188	346772	20.000	ng	0.00
76) Chrysene-d12	21.926	240	333321	20.000	ng	0.00
86) Perylene-d12	25.345	264	352700	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	172409	58.480	ng	0.00
7) Phenol-d6	7.395	99	159000	39.043	ng	0.00
23) Nitrobenzene-d5	9.440	82	310497	92.679	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	237137	158.862	ng	0.00
45) 2-Fluorobiphenyl	13.518	172	760478	74.979	ng	0.00
79) Terphenyl-d14	20.216	244	738970	44.530	ng	0.00

Target Compounds	Qvalue

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

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