

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110922\
 Data File : BG055530.D
 Acq On : 9 Nov 2022 19:29
 Operator : CG/JU
 Sample : N5360-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221171-COMP-06-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:31:26 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	8.273	152	60028	20.000	ng	0.00
21) Naphthalene-d8	11.099	136	241356	20.000	ng	# 0.00
39) Acenaphthene-d10	14.888	164	162128	20.000	ng	0.00
64) Phenanthrene-d10	17.626	188	366969	20.000	ng	0.00
76) Chrysene-d12	21.921	240	344963	20.000	ng	0.00
86) Perylene-d12	25.341	264	361578	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.799	112	190724	60.518	ng	0.00
7) Phenol-d6	7.397	99	192586	44.239	ng	0.00
23) Nitrobenzene-d5	9.442	82	317527	89.187	ng	0.00
42) 2,4,6-Tribromophenol	16.363	330	216689	136.981	ng	0.00
45) 2-Fluorobiphenyl	13.519	172	750099	69.787	ng	0.00
79) Terphenyl-d14	20.211	244	645211	37.568	ng	0.00

Target Compounds	Qvalue

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

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