

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
 Data File : BG055592.D
 Acq On : 12 Nov 2022 16:36
 Operator : CG/JU
 Sample : N5431-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221177-COMP-12-MODIFIED-ELUTRIATE

Quant Time: Nov 14 03:28:14 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	8.270	152	64518	20.000	ng	0.00
21) Naphthalene-d8	11.097	136	262784	20.000	ng	0.00
39) Acenaphthene-d10	14.886	164	197219	20.000	ng	0.00
64) Phenanthrene-d10	17.630	188	480247	20.000	ng	0.00
76) Chrysene-d12	21.925	240	477229	20.000	ng	0.00
86) Perylene-d12	25.345	264	504497	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	188165	55.551	ng	0.00
7) Phenol-d6	7.401	99	171989	36.758	ng	0.00
23) Nitrobenzene-d5	9.440	82	315618	81.422	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	290683	151.061	ng	0.00
45) 2-Fluorobiphenyl	13.517	172	777268	59.448	ng	0.00
79) Terphenyl-d14	20.215	244	1455471	61.258	ng	0.00

Target Compounds	Qvalue

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

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