

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011022\
 Data File : BF126569.D
 Acq On : 10 Jan 2022 16:36
 Operator : JU/CG
 Sample : N1054-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220020-CAPE-2-MODIFIED-ELUTRIATE

Quant Time: Jan 10 17:18:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 10 14:50:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	83796	20.000	ng	0.00
21) Naphthalene-d8	8.051	136	350742	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	164963	20.000	ng	0.00
64) Phenanthrene-d10	11.292	188	255369	20.000	ng	0.00
76) Chrysene-d12	13.939	240	147090	20.000	ng	# 0.00
86) Perylene-d12	15.398	264	156242	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.386	112	465920	79.847	ng	0.00
7) Phenol-d6	6.404	99	489910	64.822	ng	0.00
23) Nitrobenzene-d5	7.333	82	668620	97.881	ng	0.00
42) 2,4,6-Tribromophenol	10.598	330	198483	155.704	ng	0.00
45) 2-Fluorobiphenyl	9.133	172	920434	98.186	ng	0.00
79) Terphenyl-d14	12.886	244	614088	81.116	ng	0.00

Target Compounds	Qvalue

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

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