

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
 Data File : BF131101.D
 Acq On : 09 Nov 2022 22:36
 Operator : CG\JU
 Sample : N5360-18
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DISSOLVED-20221175-COMP-10-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:34:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	89864	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	347845	20.000	ng	0.00
39) Acenaphthene-d10	9.934	164	192190	20.000	ng	0.00
64) Phenanthrene-d10	11.428	188	360270	20.000	ng	0.00
76) Chrysene-d12	14.069	240	233515	20.000	ng	#-0.01
86) Perylene-d12	15.563	264	171953	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	279093	52.816	ng	0.00
7) Phenol-d6	6.516	99	298565	43.218	ng	0.00
23) Nitrobenzene-d5	7.457	82	519467	67.380	ng	0.00
42) 2,4,6-Tribromophenol	10.728	330	194895	98.123	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	887229	64.444	ng	0.00
79) Terphenyl-d14	13.010	244	461382	34.642	ng	0.00

Target Compounds	Qvalue

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

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