

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
 Data File : BF131105.D
 Acq On : 10 Nov 2022 00:35
 Operator : CG\JU
 Sample : N5360-17
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DISSOLVED-20221172-COMP-07-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:35:07 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	84054	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	317017	20.000	ng	0.00
39) Acenaphthene-d10	9.934	164	168519	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	308363	20.000	ng	0.00
76) Chrysene-d12	14.069	240	209639	20.000	ng	#-0.01
86) Perylene-d12	15.563	264	195967	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	246627	49.899	ng	0.00
7) Phenol-d6	6.510	99	243152	37.629	ng	-0.01
23) Nitrobenzene-d5	7.457	82	450069	64.055	ng	0.00
42) 2,4,6-Tribromophenol	10.728	330	181791	104.382	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	803646	66.572	ng	0.00
79) Terphenyl-d14	13.010	244	394086	32.960	ng	0.00

Target Compounds Qvalue

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

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