

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012422\
 Data File : BF126643.D
 Acq On : 24 Jan 2022 17:10
 Operator : CG/JU
 Sample : N1210-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Time: Jan 25 03:13:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 19 07:04:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	108085	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	415538	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	229668	20.000	ng	0.00
64) Phenanthrene-d10	11.498	188	397751	20.000	ng	0.00
76) Chrysene-d12	14.145	240	283392	20.000	ng	# 0.00
86) Perylene-d12	15.674	264	230278	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.575	112	439204	53.472	ng	0.00
7) Phenol-d6	6.569	99	403542	35.361	ng	-0.01
23) Nitrobenzene-d5	7.522	82	936389	88.202	ng	-0.01
42) 2,4,6-Tribromophenol	10.792	330	234625	109.162	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	1197371	79.844	ng	0.00
79) Terphenyl-d14	13.086	244	1192385	70.572	ng	0.00

Target Compounds Qvalue

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

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