

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
 Data File : BF126990.D
 Acq On : 21 Feb 2022 23:09
 Operator : CG\JU
 Sample : N1602-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220077A-BERKELEY-4A-COMP

Quant Time: Feb 22 03:53:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 21 13:16:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	90936	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	366592	20.000	ng	# 0.00
39) Acenaphthene-d10	9.957	164	191546	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	335483	20.000	ng	# 0.00
76) Chrysene-d12	14.086	240	144659	20.000	ng	#-0.01
86) Perylene-d12	15.586	264	166821	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.539	112	600179	102.008	ng	0.00
7) Phenol-d6	6.539	99	790718	99.598	ng	0.00
23) Nitrobenzene-d5	7.481	82	554896	68.323	ng	-0.01
42) 2,4,6-Tribromophenol	10.745	330	170904	77.494	ng	0.00
45) 2-Fluorobiphenyl	9.275	172	871842	67.007	ng	0.00
79) Terphenyl-d14	13.027	244	760132	77.245	ng	-0.01

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
Data File : BF126990.D
Acq On : 21 Feb 2022 23:09
Operator : CG\JU
Sample : N1602-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20220077A-BERKELEY-4A-COMP

Quant Time: Feb 22 03:53:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
Response via : Initial Calibration

