

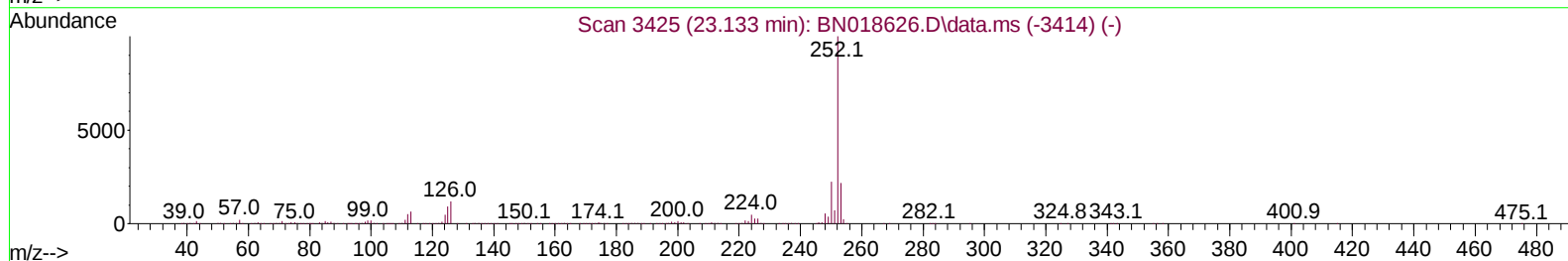
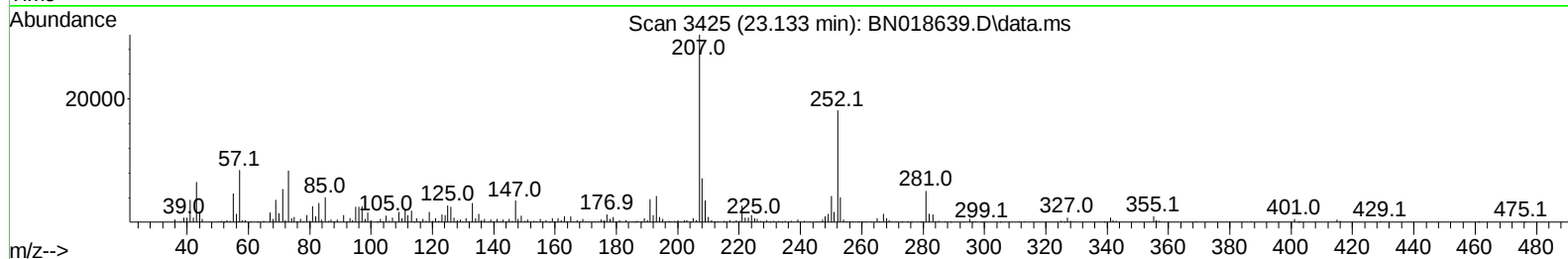
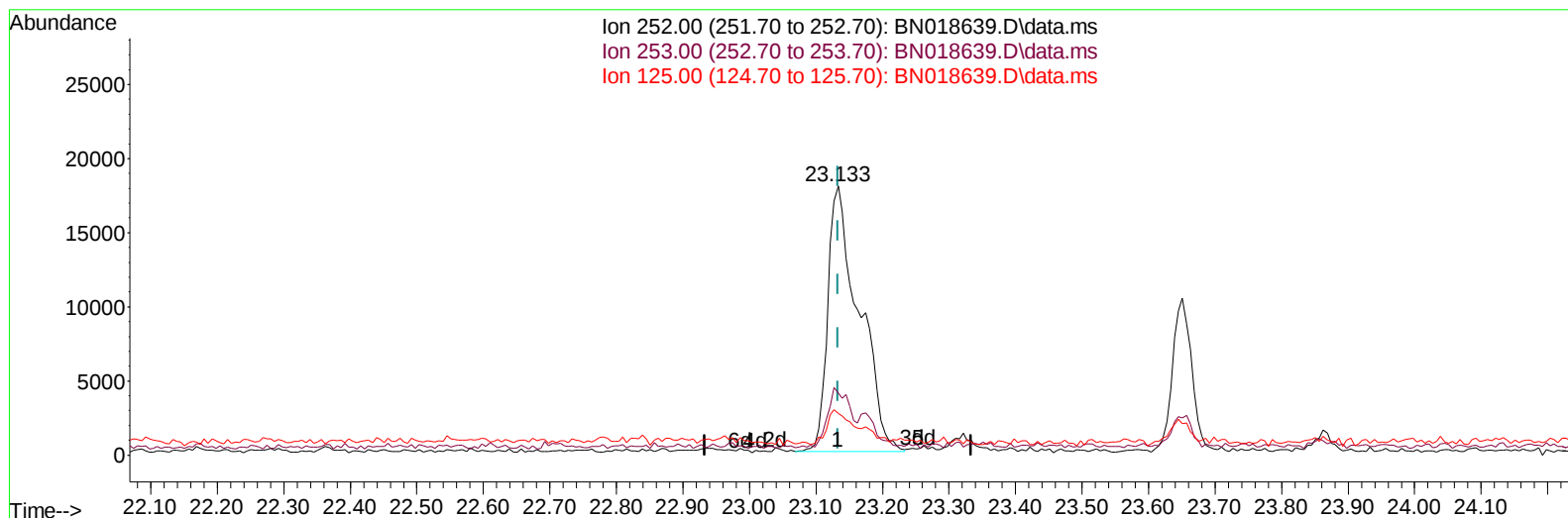
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
 Data File : BN018639.D
 Acq On : 17 Feb 2022 22:04
 Operator : CG/JU
 Sample : N1507-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGZF5

Manual IntegrationsAPPROVED

Quant Time: Feb 18 00:35:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 16 16:48:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/18/2022
 Supervised By :mohammad ahmed 02/18/2022



TIC: BN018639.D\data.ms

(90) Benzo(b)fluoranthene

23.133min (-0.000) 1.41 ng/u1

response 58192

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.20	22.96
125.00	9.00	15.55#
0.00	0.00	0.00

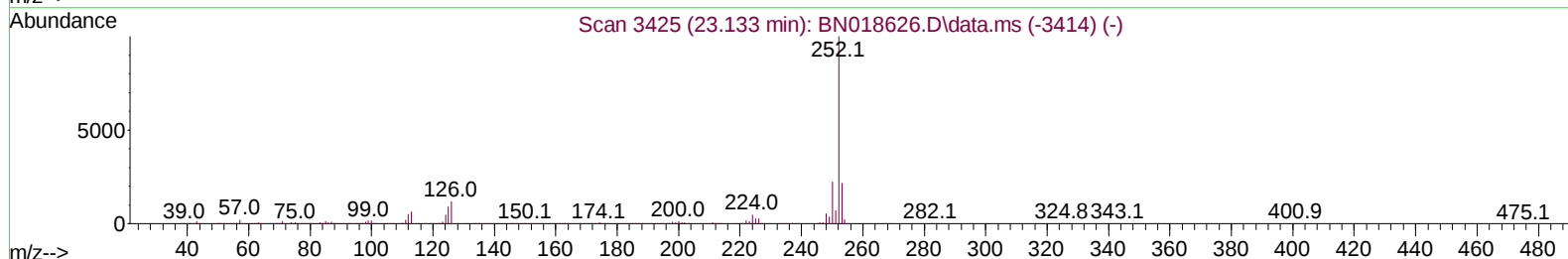
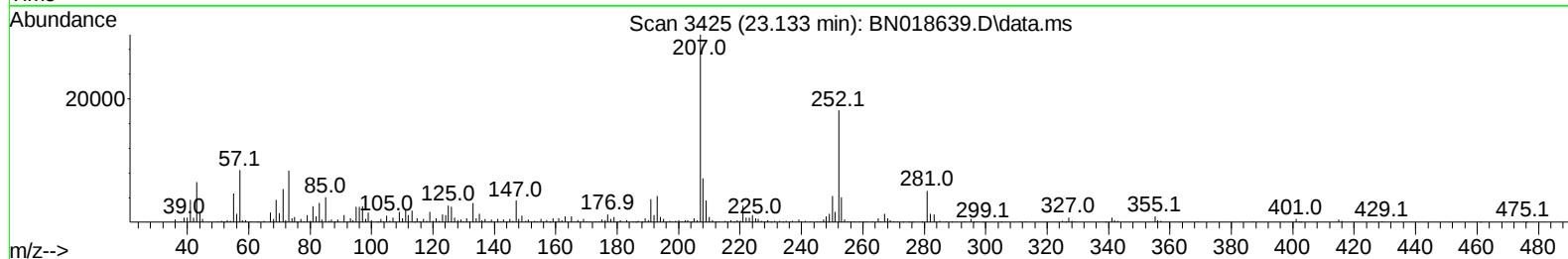
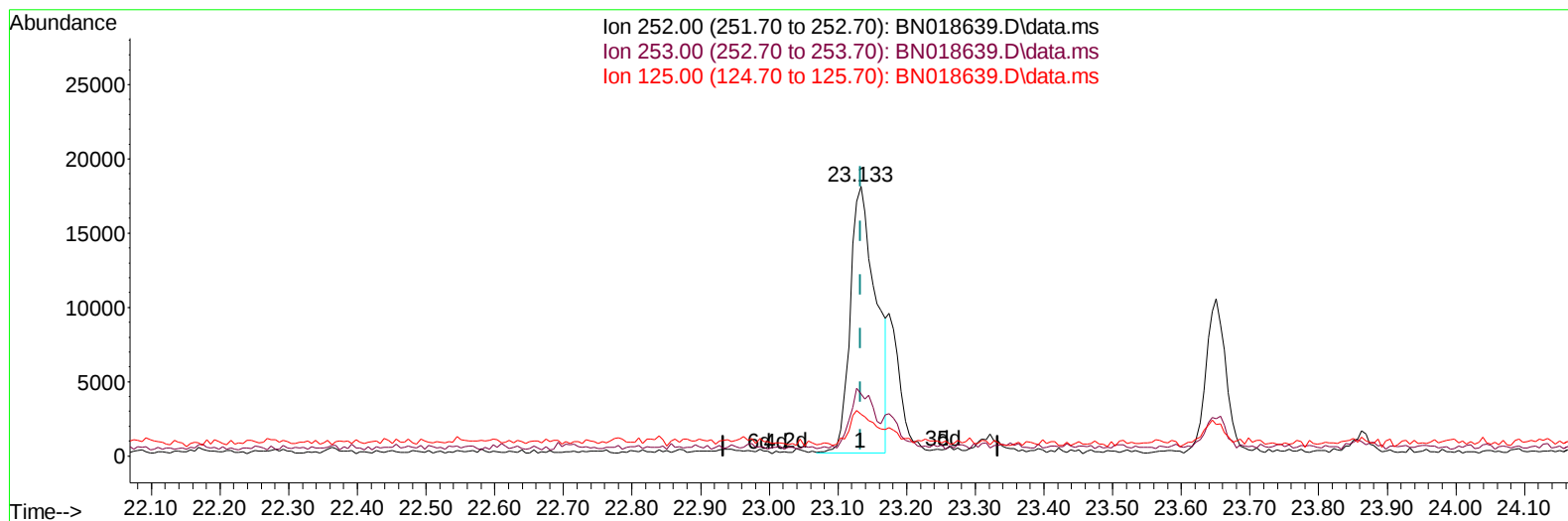
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(90) Benzo(b)fluoranthene

23.133min (-0.000) 1.13 ng/ul m

response 46609

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.20	22.96
125.00	9.00	15.55#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.916	152		214429	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136		832489	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164		467174	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188		823135	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240		655533	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264		612309	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	0.000	96		0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	0.000	99		0	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67		0	0.000	ng/ul	
11) 2-Chlorophenol-d4	0.000	132		0	0.000	ng/ul	
15) 4-Methylphenol-d8	0.000	113		0	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128		0	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143		0	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165		0	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131		0d	0.000	ng/ul	
46) Dimethylphthalate-d6	0.000	166		0	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160		0	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143		0	0.000	ng/ul	
60) Fluorene-d10	0.000	176		0	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200		0	0.000	ng/ul	
73) Anthracene-d10	0.000	188		0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212		0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264		0d	0.000	ng/ul	
Target Compounds							
90) Benzo(b)fluoranthene	23.133	252		46609m	1.129	ng/ul	Qvalue

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

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