

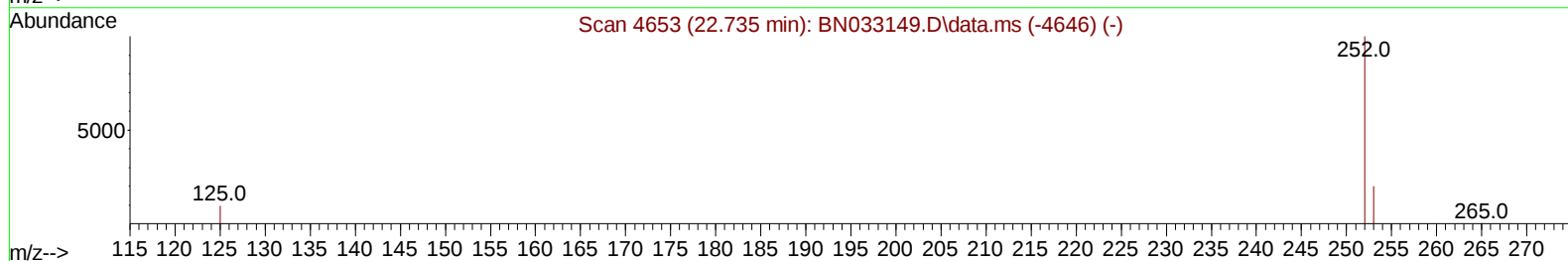
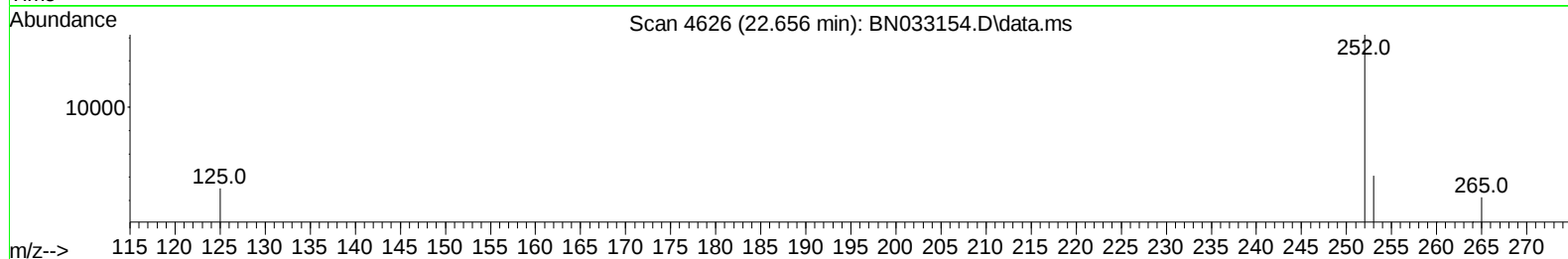
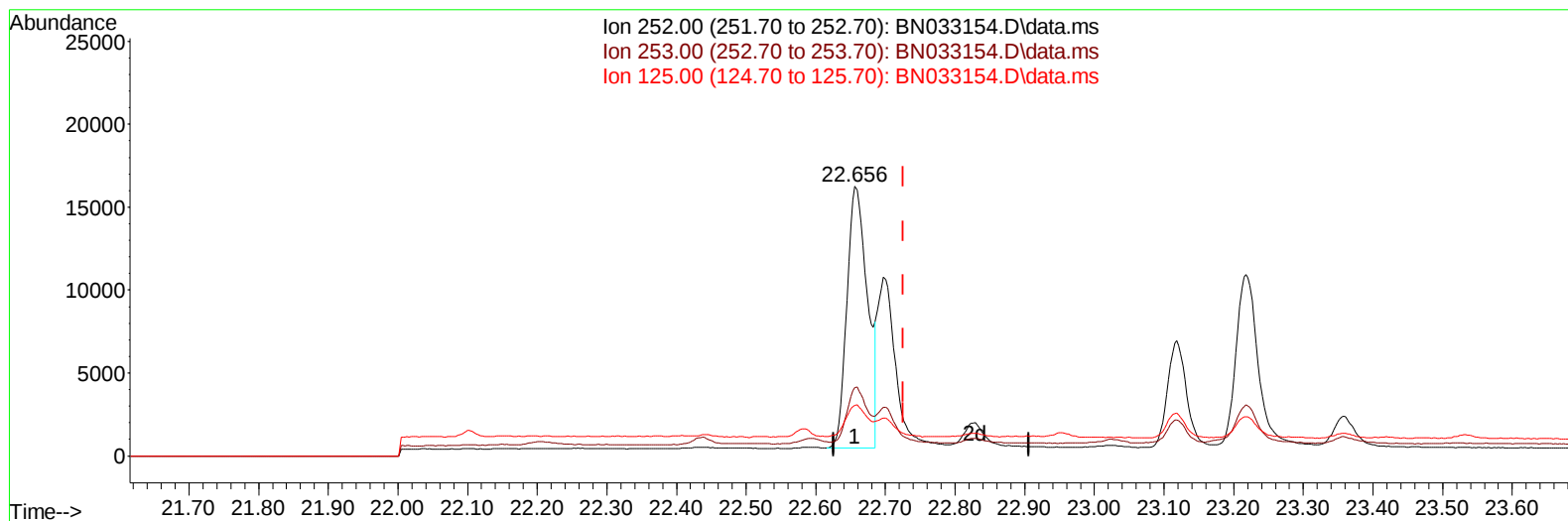
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
Data File : BN033154.D
Acq On : 01 Aug 2024 05:54
Operator : MA/JU
Sample : P3265-11MS
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4D42MS

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/01/2024
Supervised By :mohammad ahmed 08/13/2024

Quant Time: Aug 01 07:42:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 30 14:49:49 2024
Response via : Initial Calibration



TIC: BN033154.D\data.ms

(25) Benzo(k)fluoranthene

22.656min (-0.070) 0.97 ng/u1

response 32738

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.10	25.36
125.00	26.10	18.88#
0.00	0.00	0.00

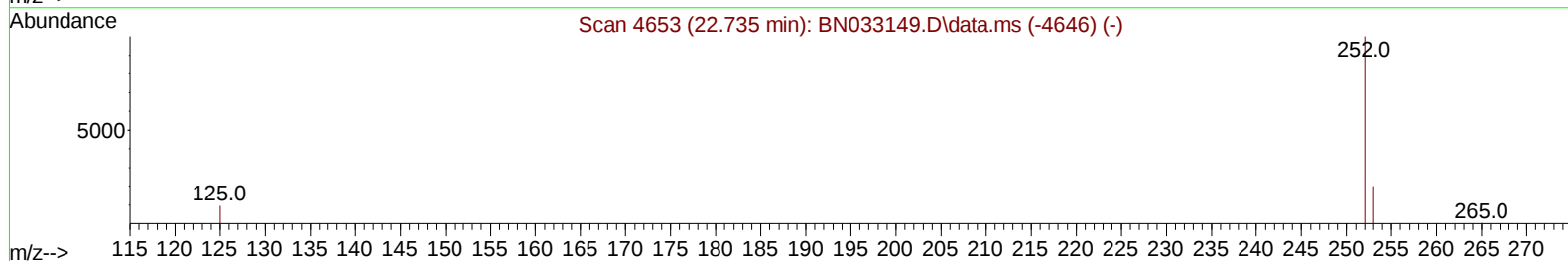
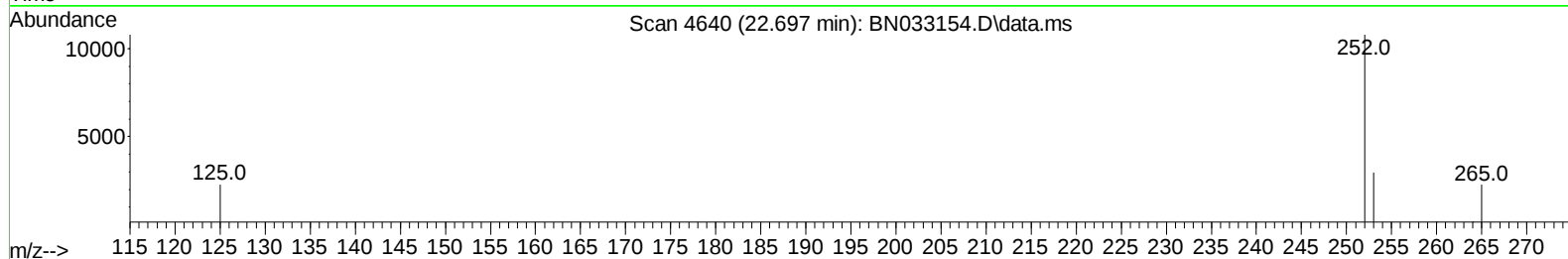
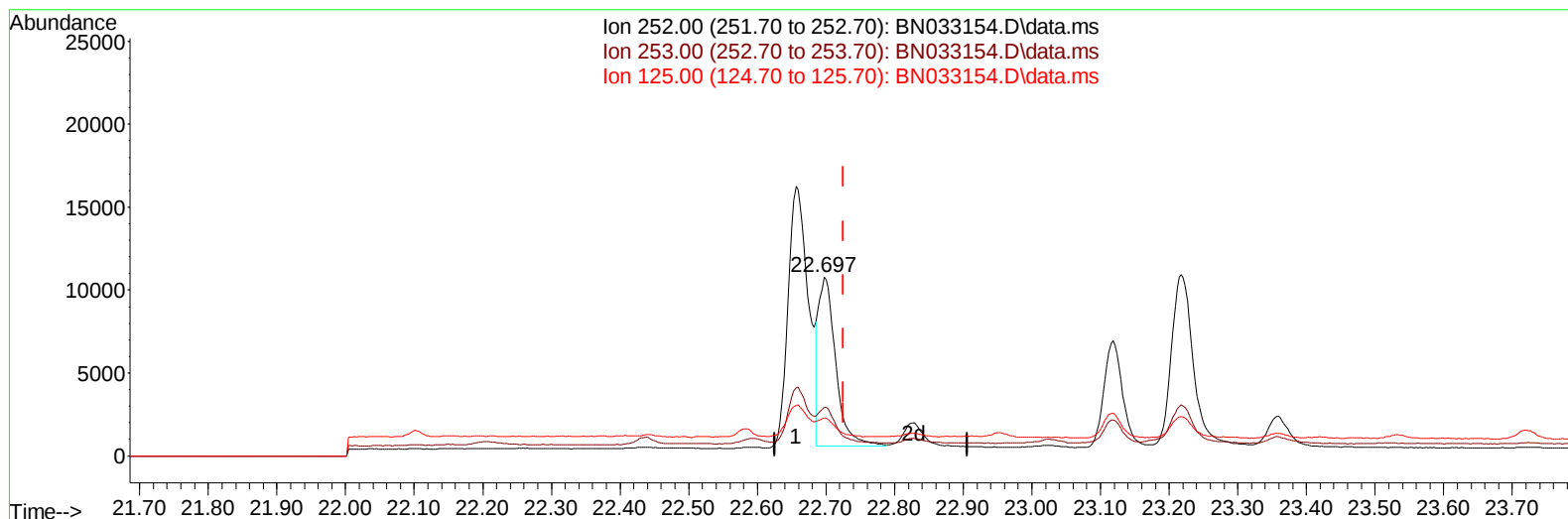
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TIC: BN033154.D\data.ms

(25) Benzo(k)fluoranthene

22.697min (-0.029) 0.52 ng/u1 m

response 17610

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.10	27.31
125.00	26.10	21.13
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	7.516	152		2331	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.279	136		7933	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.143	164		4724	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.912	188		10091	0.400	ng/ul	-0.03
17) Chrysene-d12	21.128	240		9233	0.400	ng/ul	-0.02
23) Perylene-d12	23.308	264		9640	0.400	ng/ul	-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.056	96		1185	0.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.868	152		3211	0.248	ng/ul	-0.06
18) Fluoranthene-d10	18.949	212		8264	0.316	ng/ul	-0.02
Target Compounds							
							Qvalue
2) 1,4-Dioxane	3.086	88		3512	1.045	ng/ul#	73
5) Naphthalene	10.328	128		6316	0.310	ng/ul	98
7) 2-Methylnaphthalene	11.950	142		3701	0.268	ng/ul	98
8) 1-Methylnaphthalene	12.165	142		3829	0.257	ng/ul	100
10) Acenaphthylene	13.870	152		6433	0.333	ng/ul	97
11) Acenaphthene	14.203	153		5207	0.358	ng/ul	99
12) Fluorene	15.217	166		5586	0.330	ng/ul	99
14) Pentachlorophenol	16.603	266		1021	0.497	ng/ul	99
15) Phenanthrene	16.954	178		23060	0.852	ng/ul	98
16) Anthracene	17.051	178		9642	0.396	ng/ul	99
19) Fluoranthene	18.977	202		45818	1.507	ng/ul	98
20) Pyrene	19.349	202		43738	1.342	ng/ul	96
21) Benzo(a)anthracene	21.113	228		25719	0.817	ng/ul	98
22) Chrysene	21.166	228		27413	0.746	ng/ul	100
24) Benzo(b)fluoranthene	22.656	252		32738	1.163	ng/ul	89
25) Benzo(k)fluoranthene	22.697	252		17610m	0.522	ng/ul	
26) Benzo(a)pyrene	23.218	252		22323	0.842	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	25.500	276		20666	0.513	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.494	278		9971	0.312	ng/ul	99
29) Benzo(g,h,i)perylene	26.174	276		18758	0.586	ng/ul	96

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

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