

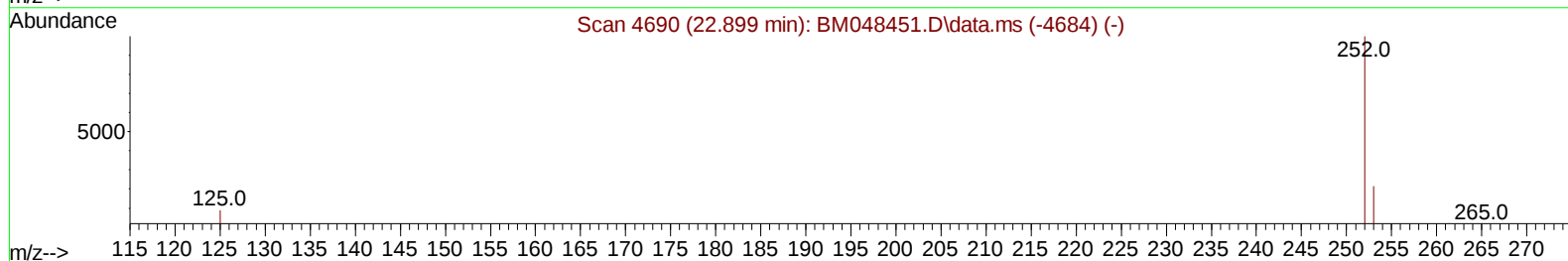
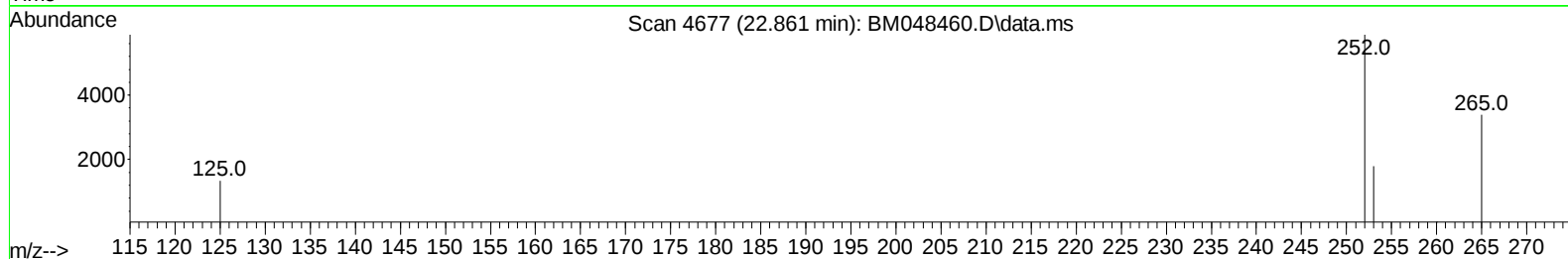
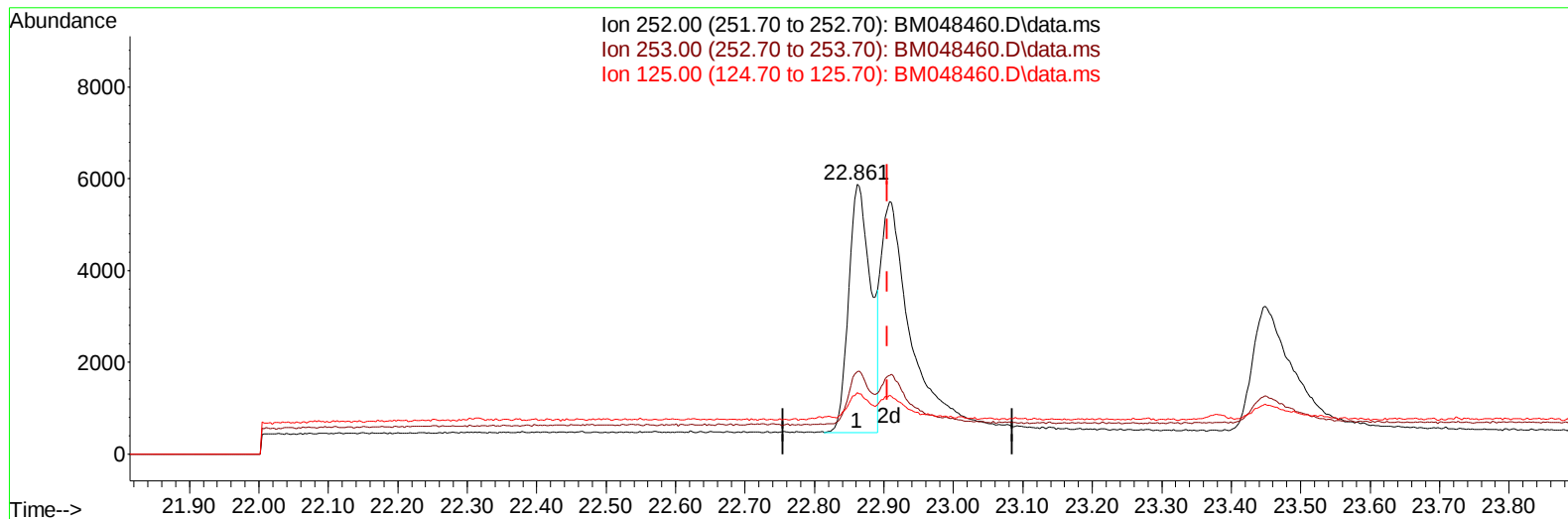
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110424\
Data File : BM048460.D
Acq On : 05 Nov 2024 16:31
Operator : RC/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC0.4

Manual IntegrationsAPPROVED

Quant Time: Nov 05 17:48:16 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 05 05:12:02 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/06/2024
Supervised By :mohammad ahmed 11/06/2024



TIC: BM048460.D\data.ms

(25) Benzo(k)fluoranthene

22.861min (-0.044) 0.32 ng/u1

response 11846

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	29.20	30.45
125.00	19.10	22.80
0.00	0.00	0.00

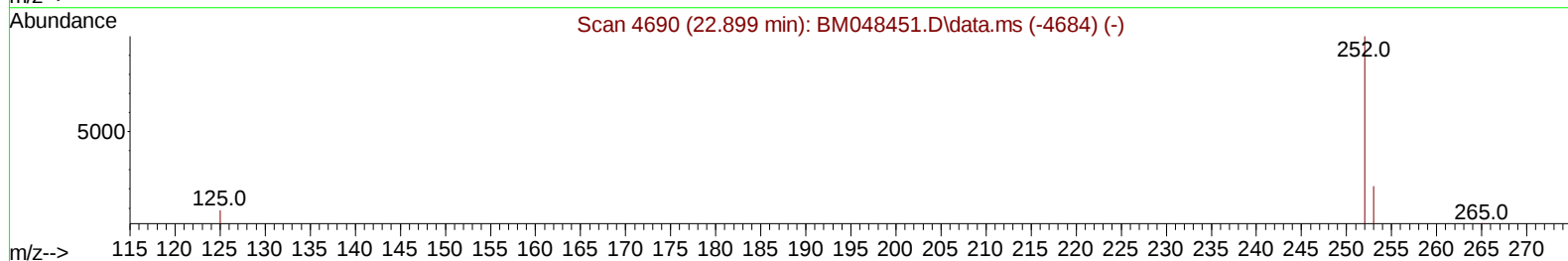
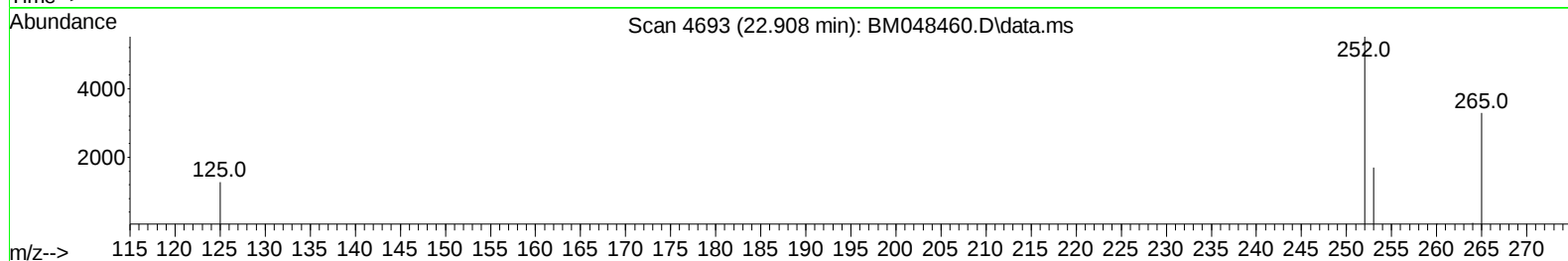
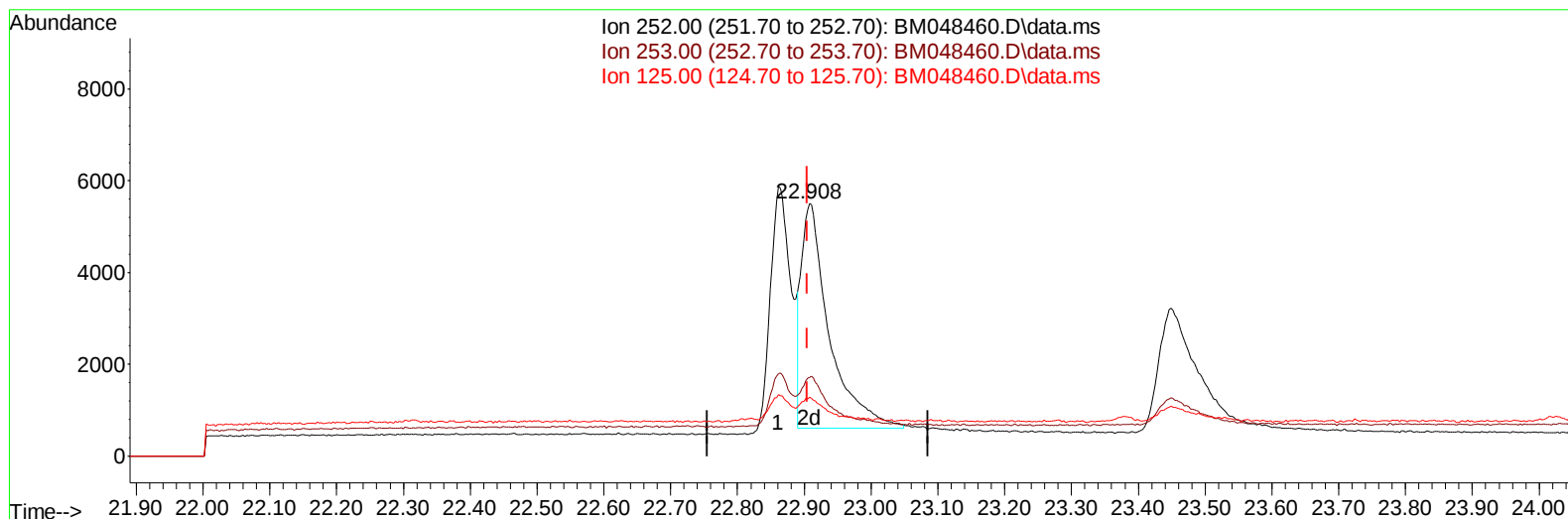
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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TIC: BM048460.D\data.ms

(25) Benzo(k)fluoranthene

22.908min (+ 0.003) 0.38 ng/ul m

response 14080

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	29.20	31.01
125.00	19.10	23.24#
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	7.758	152		2824	0.400	ng/ul	0.00
4) Naphthalene-d8	10.530	136		9796	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.359	164		5561	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.124	188		10487	0.400	ng/ul	0.00
17) Chrysene-d12	21.298	240		8548	0.400	ng/ul	# 0.00
23) Perylene-d12	23.533	264		9087	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.160	96		1727	0.498	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.152	152		4683	0.376	ng/ul	0.01
18) Fluoranthene-d10	19.142	212		9861	0.454	ng/ul	0.00
Target Compounds							
							Qvalue
2) 1,4-Dioxane	3.193	88		2031	0.540	ng/ul#	89
5) Naphthalene	10.579	128		9881	0.386	ng/ul	98
7) 2-Methylnaphthalene	12.223	142		5590	0.367	ng/ul	99
8) 1-Methylnaphthalene	12.416	142		6367	0.392	ng/ul	99
10) Acenaphthylene	14.090	152		8335	0.388	ng/ul	99
11) Acenaphthene	14.423	153		6876	0.399	ng/ul	97
12) Fluorene	15.437	166		7202	0.376	ng/ul	99
14) Pentachlorophenol	16.791	266		863	0.296	ng/ul	94
15) Phenanthrene	17.167	178		10154	0.390	ng/ul	100
16) Anthracene	17.276	178		8994	0.369	ng/ul	98
19) Fluoranthene	19.169	202		13264	0.450	ng/ul	97
20) Pyrene	19.532	202		14234	0.450	ng/ul	97
21) Benzo(a)anthracene	21.289	228		8734	0.341	ng/ul	99
22) Chrysene	21.330	228		15959	0.453	ng/ul	100
24) Benzo(b)fluoranthene	22.861	252		11843	0.381	ng/ul	96
25) Benzo(k)fluoranthene	22.908	252		14080m	0.383	ng/ul	
26) Benzo(a)pyrene	23.449	252		11368	0.390	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.832	276		15102	0.370	ng/ul	96
28) Dibenzo(a,h)anthracene	25.853	278		11720	0.369	ng/ul	93
29) Benzo(g,h,i)perylene	26.503	276		13494	0.393	ng/ul	99

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

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