

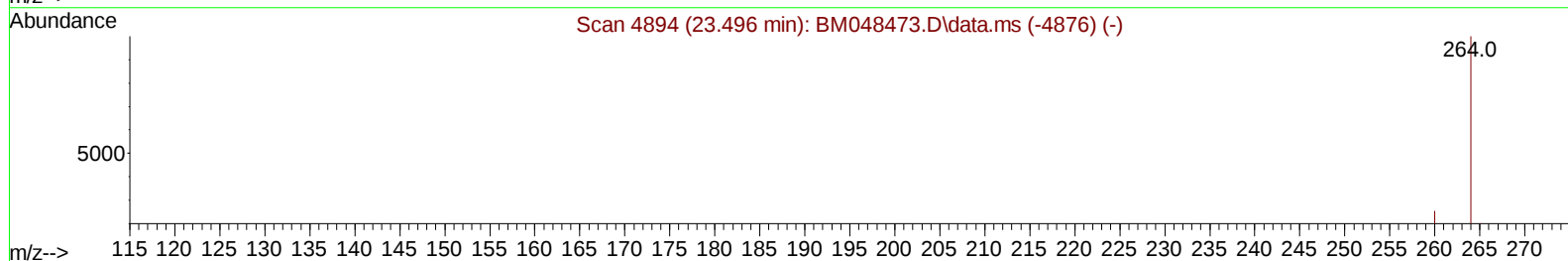
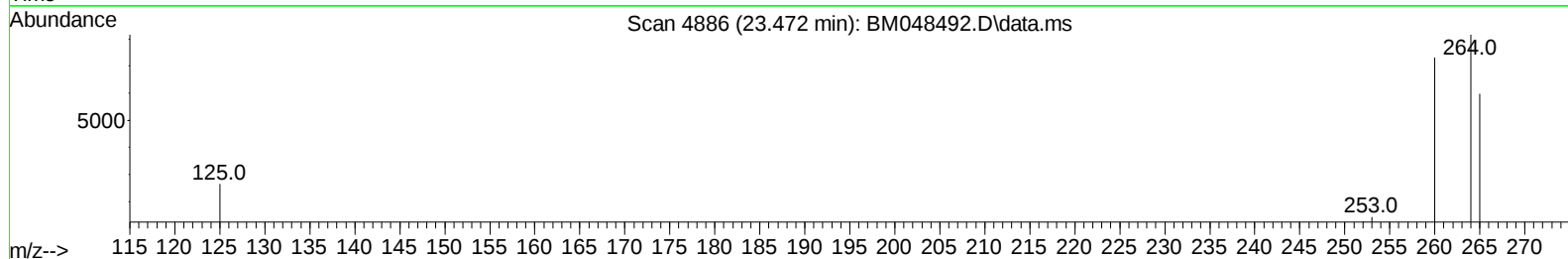
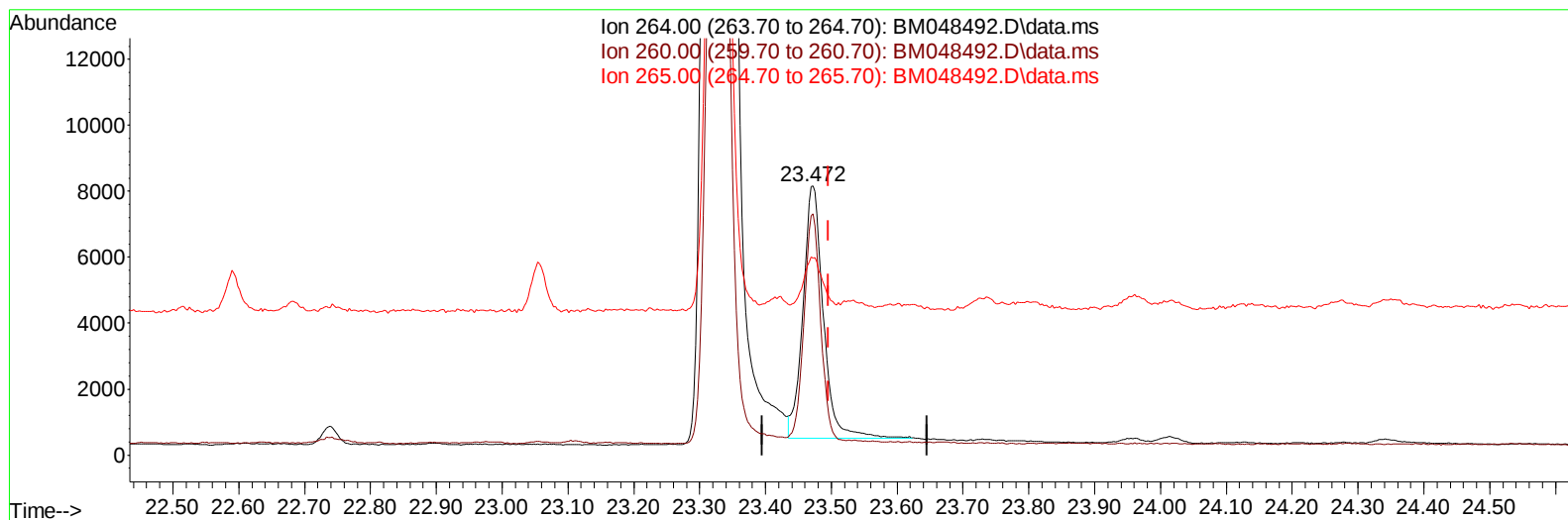
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
Data File : BM048492.D
Acq On : 07 Nov 2024 00:48
Operator : RC/JU
Sample : P4634-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CC0P9

Manual IntegrationsAPPROVED

Quant Time: Nov 07 01:20:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 06 14:39:10 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/07/2024
Supervised By :mohammad ahmed 11/11/2024



TIC: BM048492.D\data.ms

(23) Perylene-d12 (I)

23.472min (-0.023) 0.40 ng/u1

response 16329

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	26.30	89.65#
265.00	119.80	73.23#
0.00	0.00	0.00

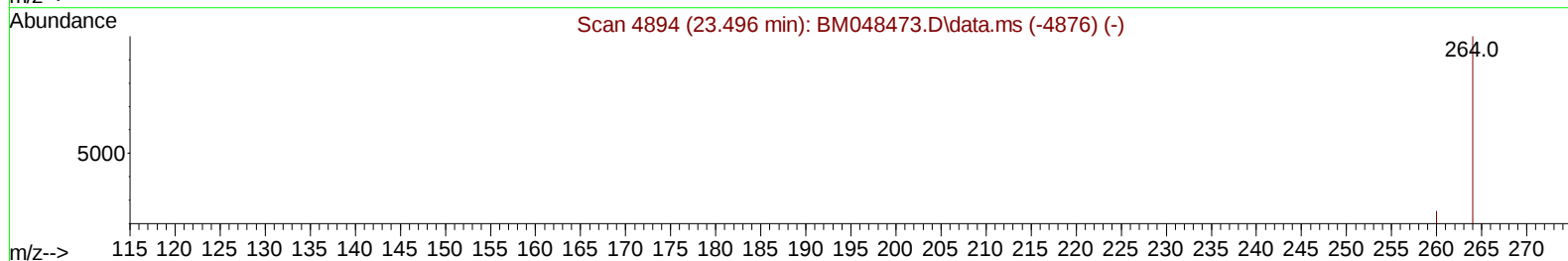
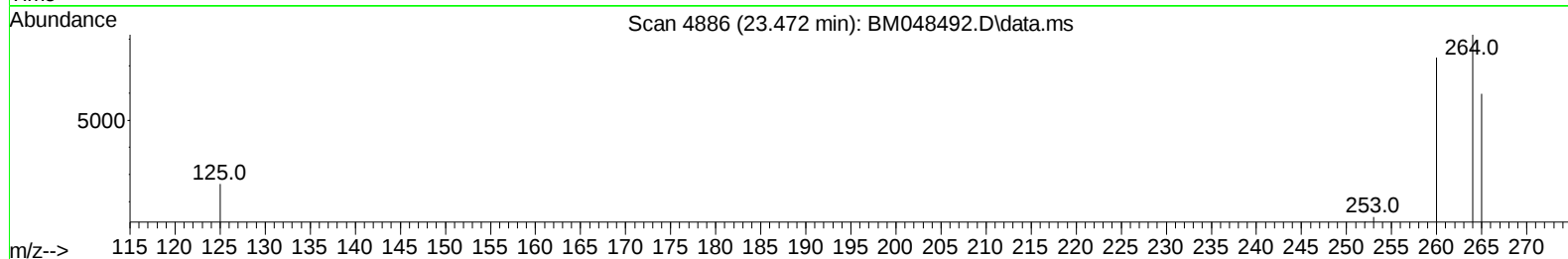
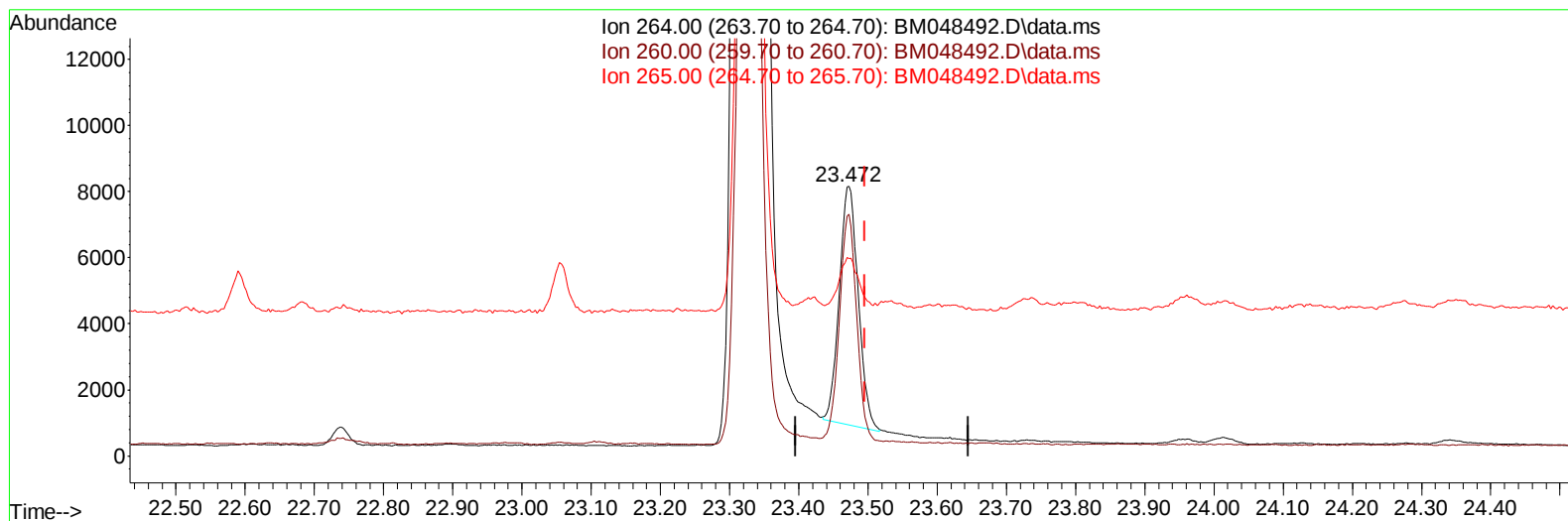
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ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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TIC: BM048492.D\data.ms

(23) Perylene-d12 (I)

23.472min (-0.023) 0.40 ng/ul m

response 13719

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	26.30	89.65#
265.00	119.80	73.23#
0.00	0.00	0.00

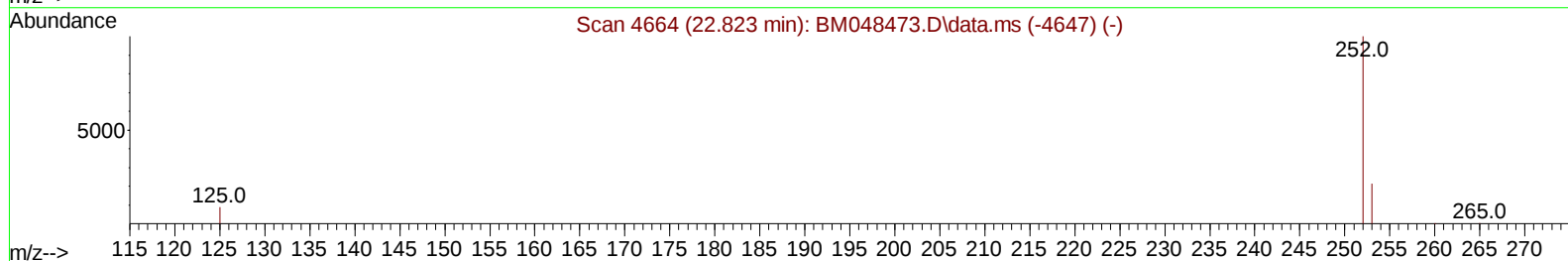
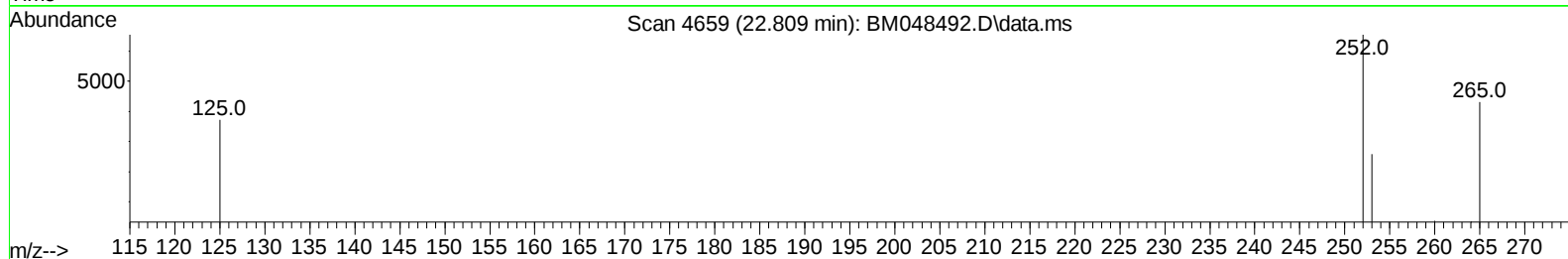
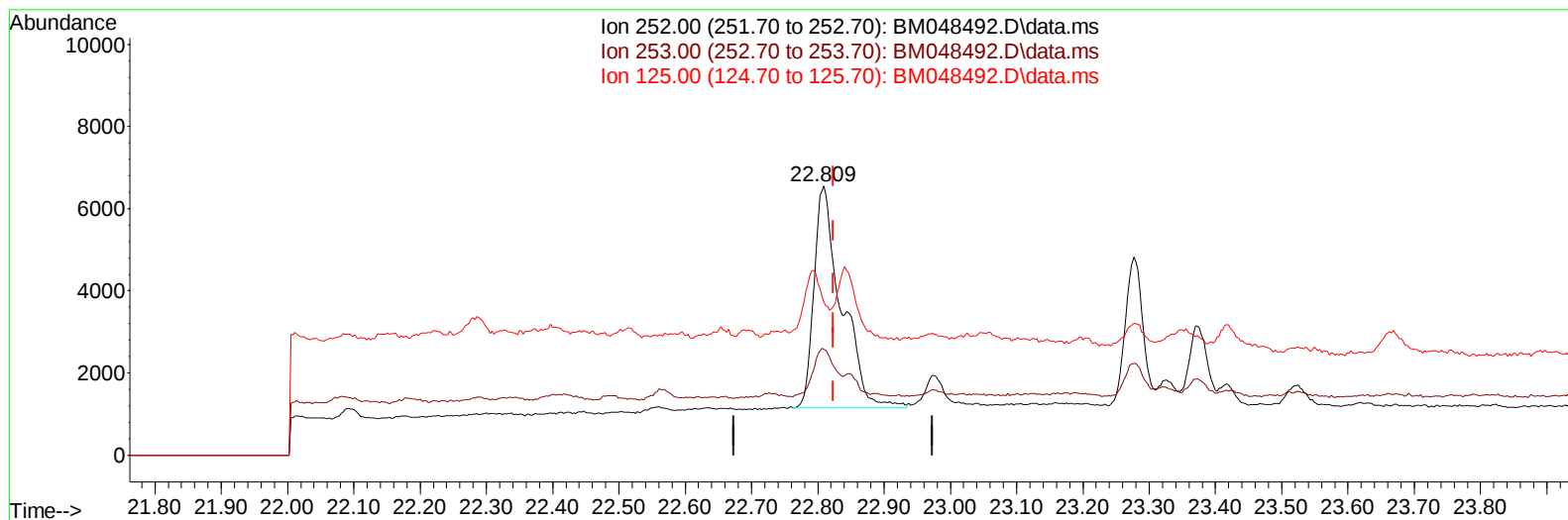
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Data File : BM048492.D
Acq On : 07 Nov 2024 00:48
Operator : RC/JU
Sample : P4634-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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TIC: BM048492.D\data.ms

(24) Benzo(b)fluoranthene

22.809min (-0.015) 0.36 ng/u1

response 15594

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	30.80	39.37
125.00	21.60	56.84#
0.00	0.00	0.00

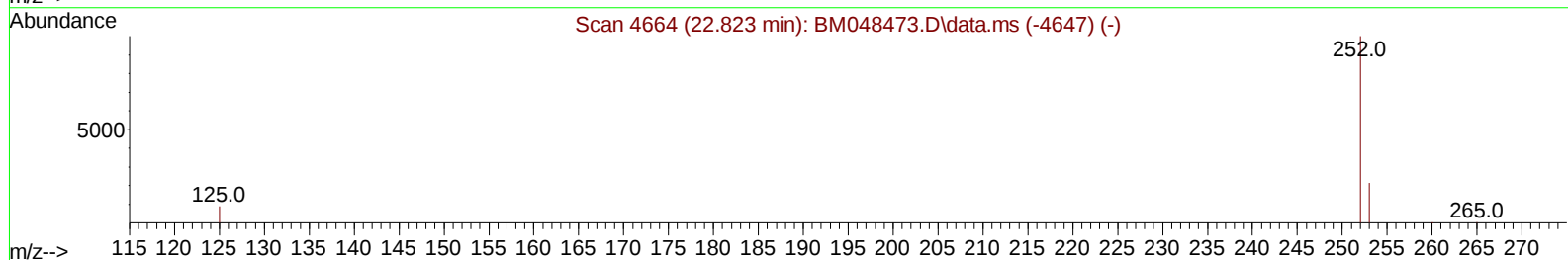
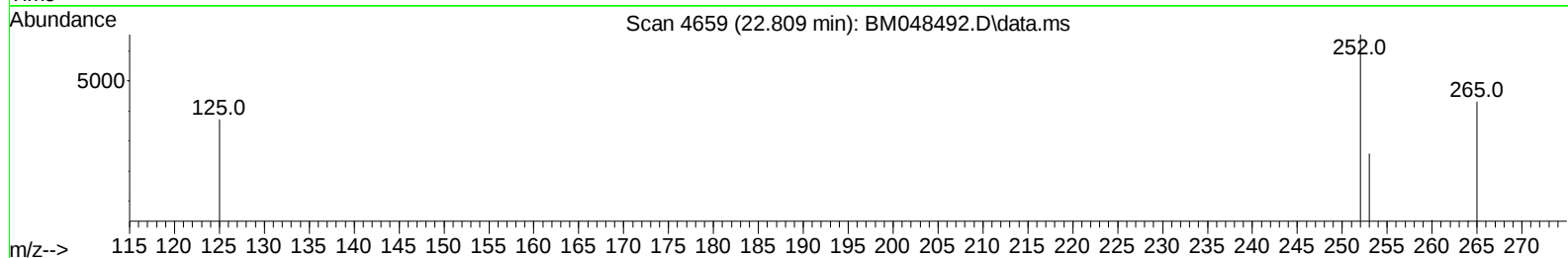
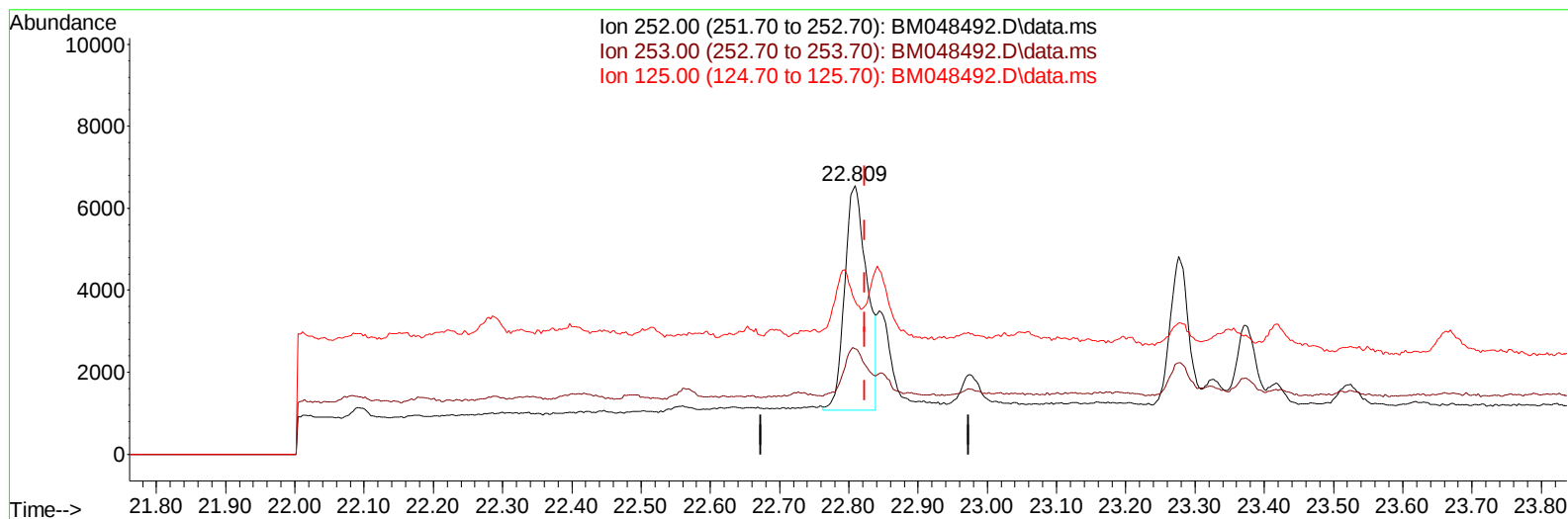
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
Data File : BM048492.D
Acq On : 07 Nov 2024 00:48
Operator : RC/JU
Sample : P4634-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CC0P9

Manual IntegrationsAPPROVED

Quant Time: Nov 07 01:20:51 2024
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TIC: BM048492.D\data.ms

(24) Benzo(b)fluoranthene

22.809min (-0.015) 0.29 ng/u1 m

response 12260

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	30.80	39.37
125.00	21.60	56.84#
0.00	0.00	0.00

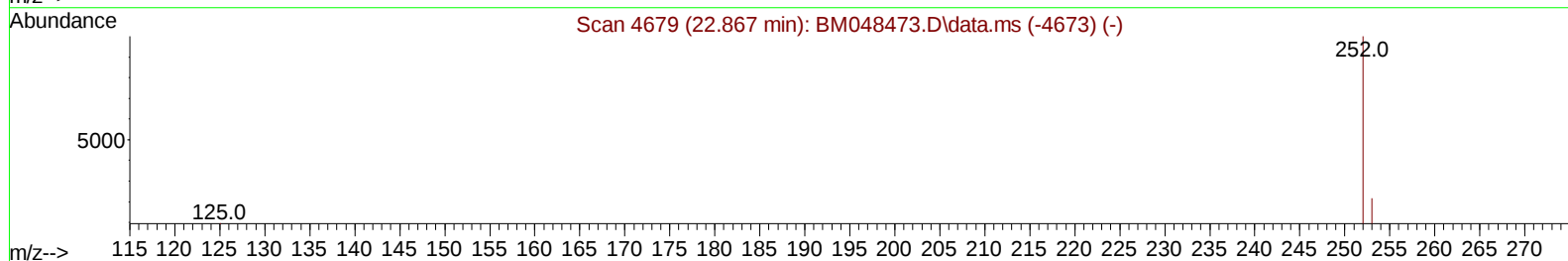
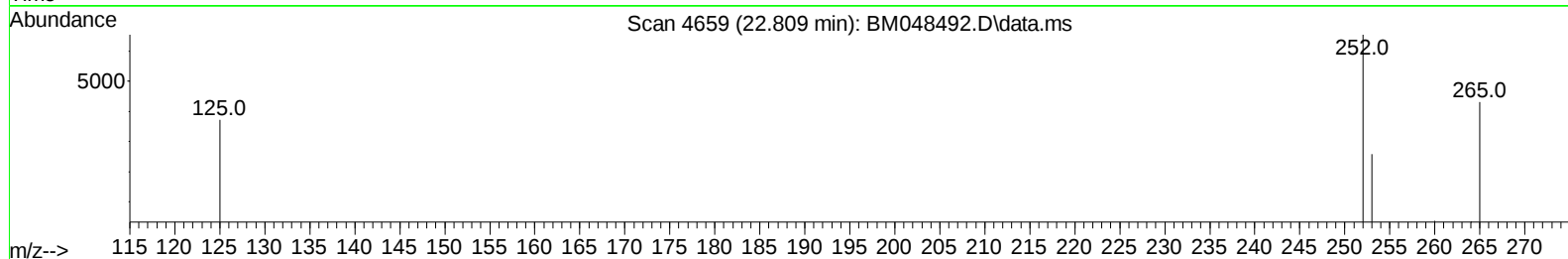
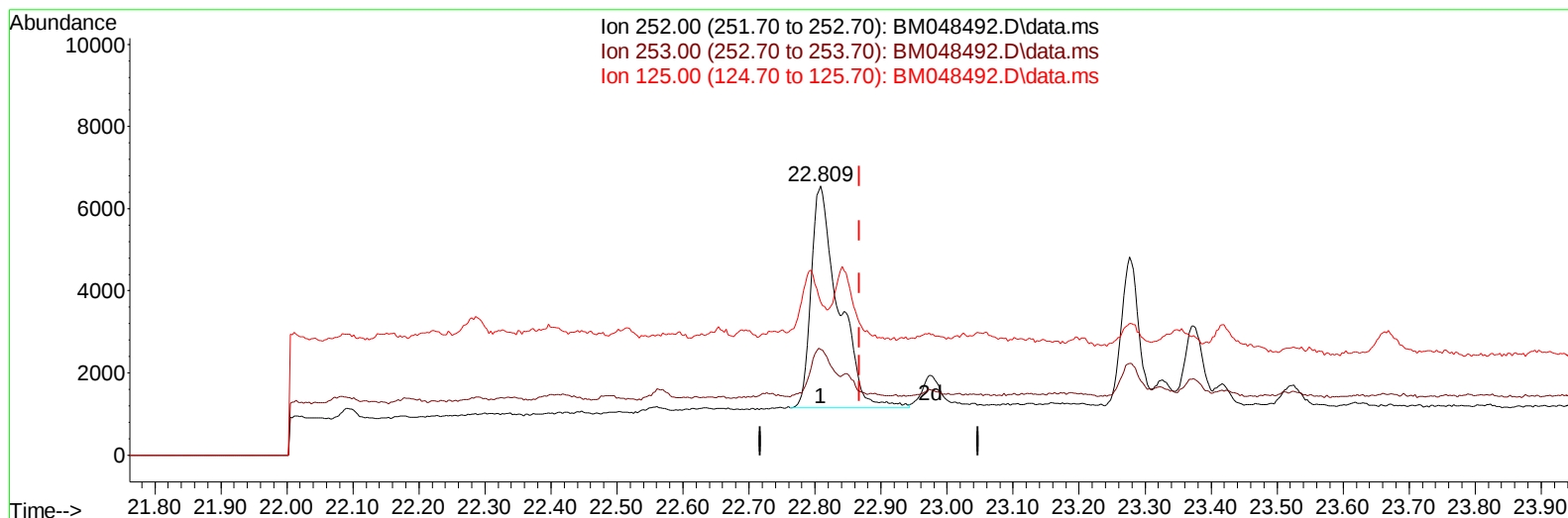
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
Data File : BM048492.D
Acq On : 07 Nov 2024 00:48
Operator : RC/JU
Sample : P4634-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CC0P9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
Supervised By :mohammad ahmed 11/11/2024

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TIC: BM048492.D\data.ms

(25) Benzo(k)fluoranthene

22.809min (-0.059) 0.31 ng/u1

response 15640

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.20	39.37#
125.00	21.70	56.84#
0.00	0.00	0.00

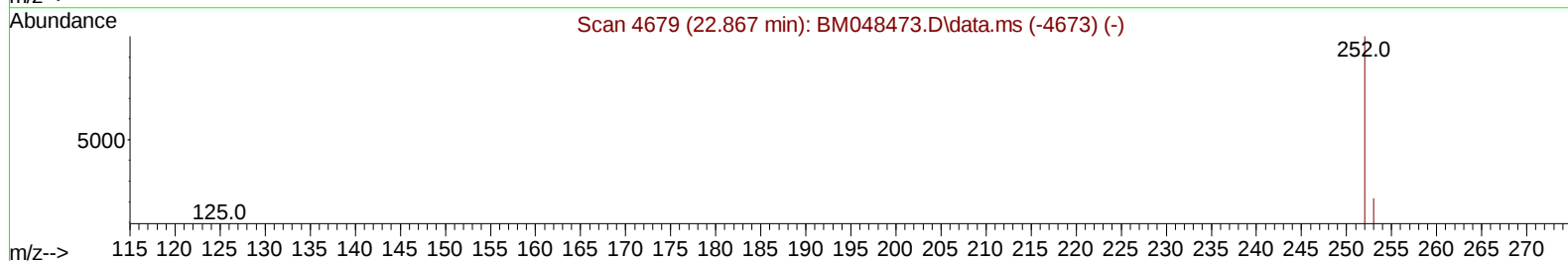
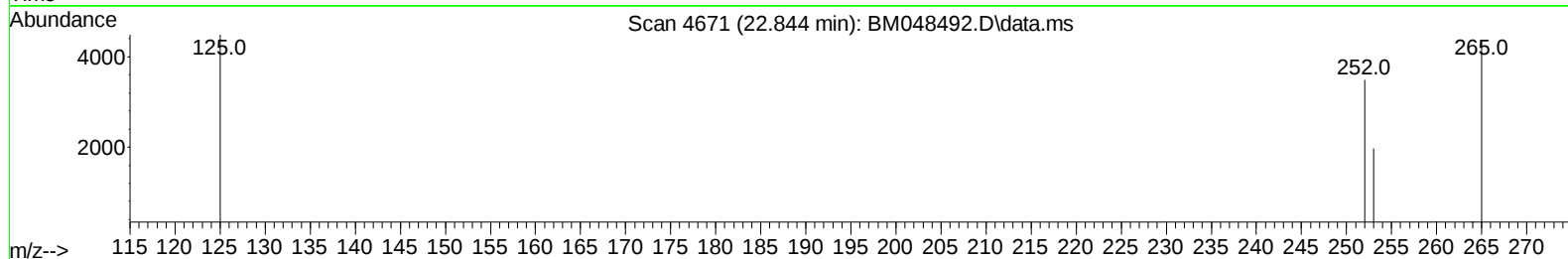
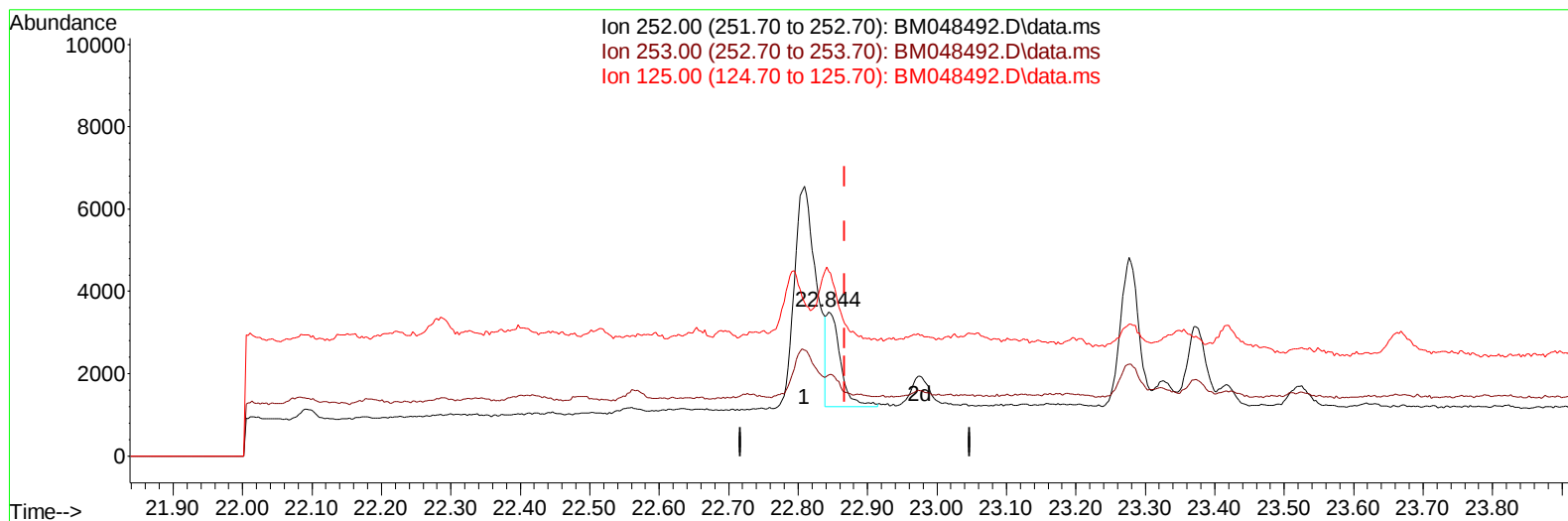
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
Data File : BM048492.D
Acq On : 07 Nov 2024 00:48
Operator : RC/JU
Sample : P4634-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CC0P9

Manual IntegrationsAPPROVED

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TIC: BM048492.D\data.ms

(25) Benzo(k)fluoranthene

22.844min (-0.023) 0.07 ng/ul m

response 3295

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.20	56.36#
125.00	21.70	128.40#
0.00	0.00	0.00

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 Sample : P4634-10
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CC0P9

Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.674	152	4766	0.400	ng/ul	0.00
4) Naphthalene-d8	10.462	136	14960	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.320	164	7965	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.064	188	15516	0.400	ng/ul	-0.02
17) Chrysene-d12	21.251	240	13248	0.400	ng/ul	#-0.01
23) Perylene-d12	23.472	264	13719m	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	9710	1.689	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.063	152	2251	0.116	ng/ul	-0.01
18) Fluoranthene-d10	19.098	212	5199	0.140	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.506	128	3146	0.084	ng/ul#	91
7) 2-Methylnaphthalene	12.134	142	849	0.036	ng/ul	88
8) 1-Methylnaphthalene	12.354	142	1020	0.042	ng/ul	97
10) Acenaphthylene	14.038	152	1033	0.034	ng/ul#	71
14) Pentachlorophenol	16.730	266	118	0.021	ng/ul	99
15) Phenanthrene	17.106	178	6525	0.160	ng/ul	96
16) Anthracene	17.199	178	1307	0.034	ng/ul#	69
19) Fluoranthene	19.126	202	7189	0.147	ng/ul#	88
20) Pyrene	19.489	202	7411	0.139	ng/ul#	90
21) Benzo(a)anthracene	21.234	228	5433	0.125	ng/ul#	77
22) Chrysene	21.286	228	10044	0.193	ng/ul#	84
24) Benzo(b)fluoranthene	22.809	252	12260m	0.287	ng/ul	
25) Benzo(k)fluoranthene	22.844	252	3295m	0.066	ng/ul	
26) Benzo(a)pyrene	23.370	252	2718	0.065	ng/ul#	29
27) Indeno(1,2,3-cd)pyrene	25.719	276	4668	0.080	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.722	278	2478	0.055	ng/ul#	1
29) Benzo(g,h,i)perylene	26.403	276	5445	0.111	ng/ul#	67

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

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