

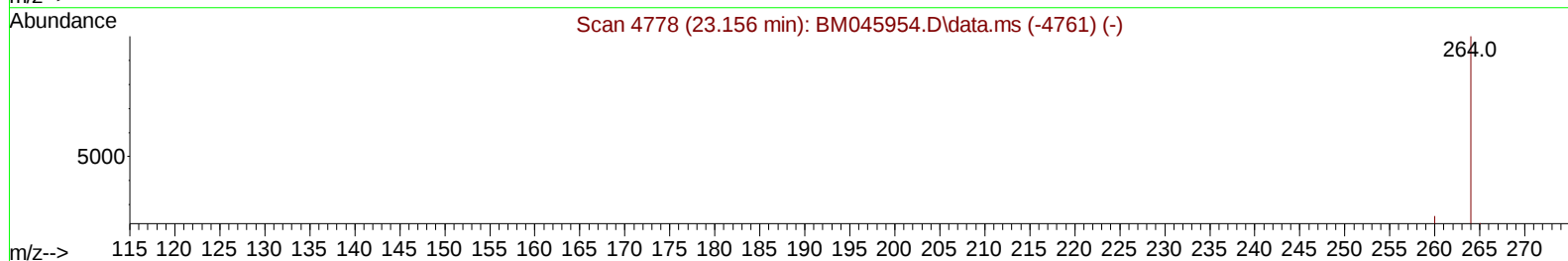
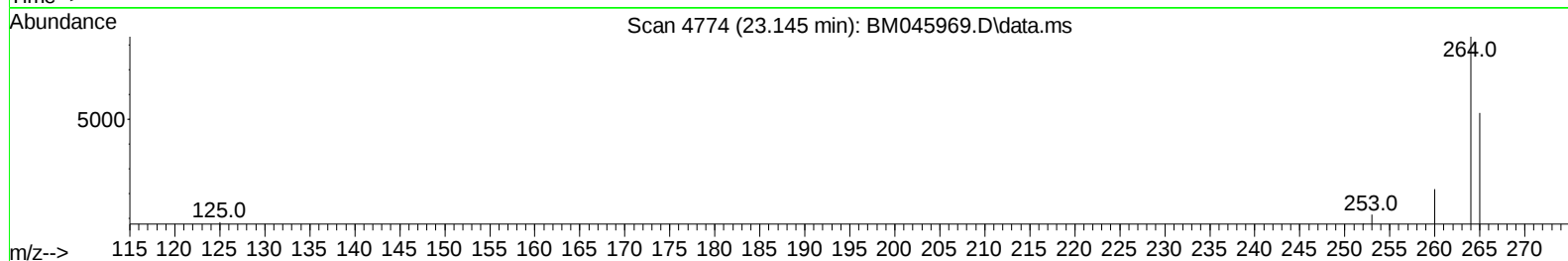
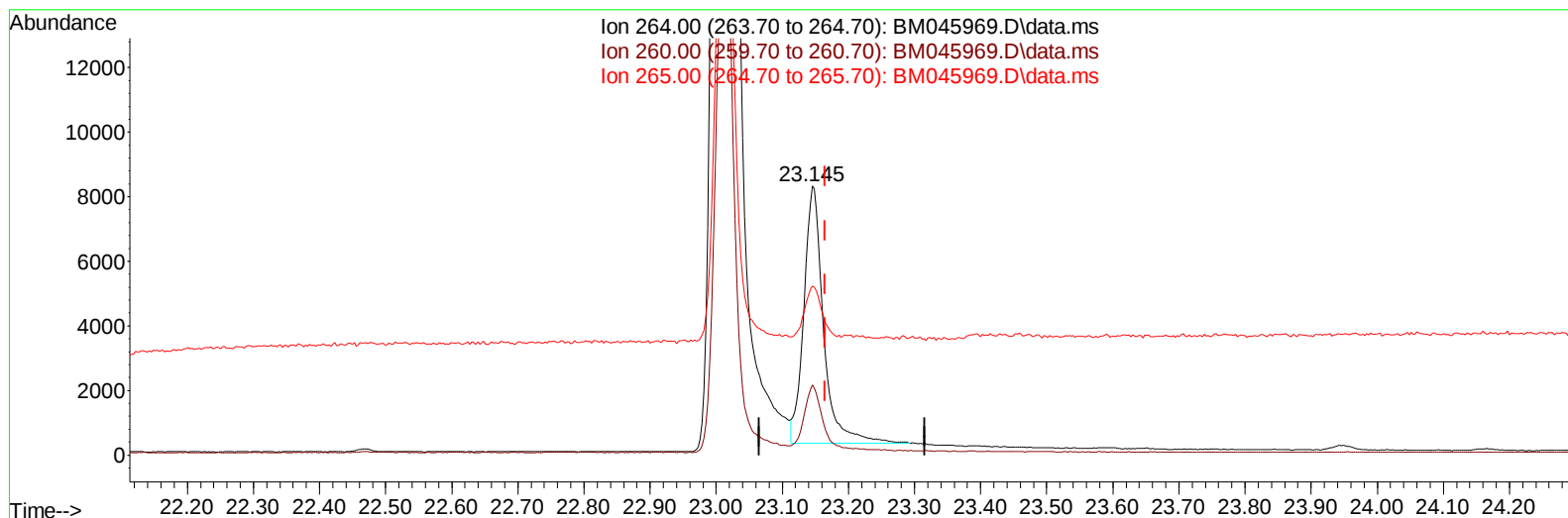
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061024\
 Data File : BM045969.D
 Acq On : 11 Jun 2024 02:12
 Operator : MA/JU
 Sample : P2642-17DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4B39DL

Manual IntegrationsAPPROVED

Quant Time: Jun 11 02:42:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM060624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 22:47:12 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/11/2024
 Supervised By :mohammad ahmed 06/13/2024



TIC: BM045969.D\data.ms

(23) Perylene-d12 (I)

23.145min (-0.020) 0.40 ng/u1

response 16298

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	25.70	26.13
265.00	76.60	62.92
0.00	0.00	0.00

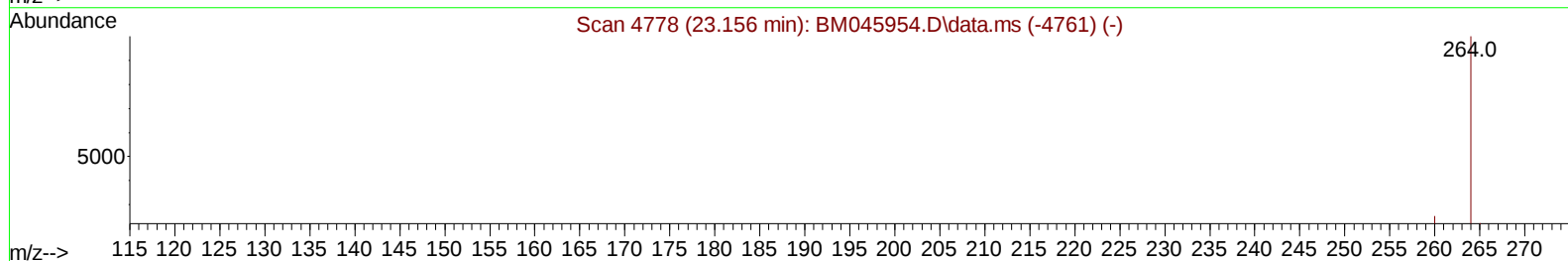
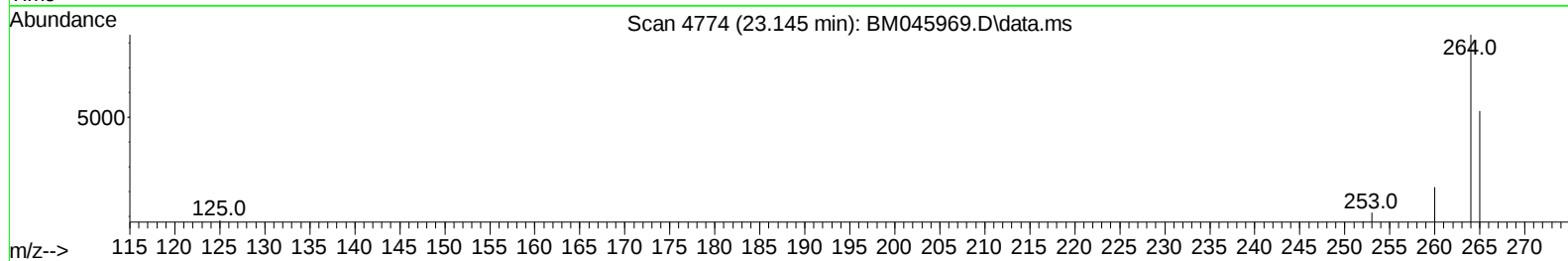
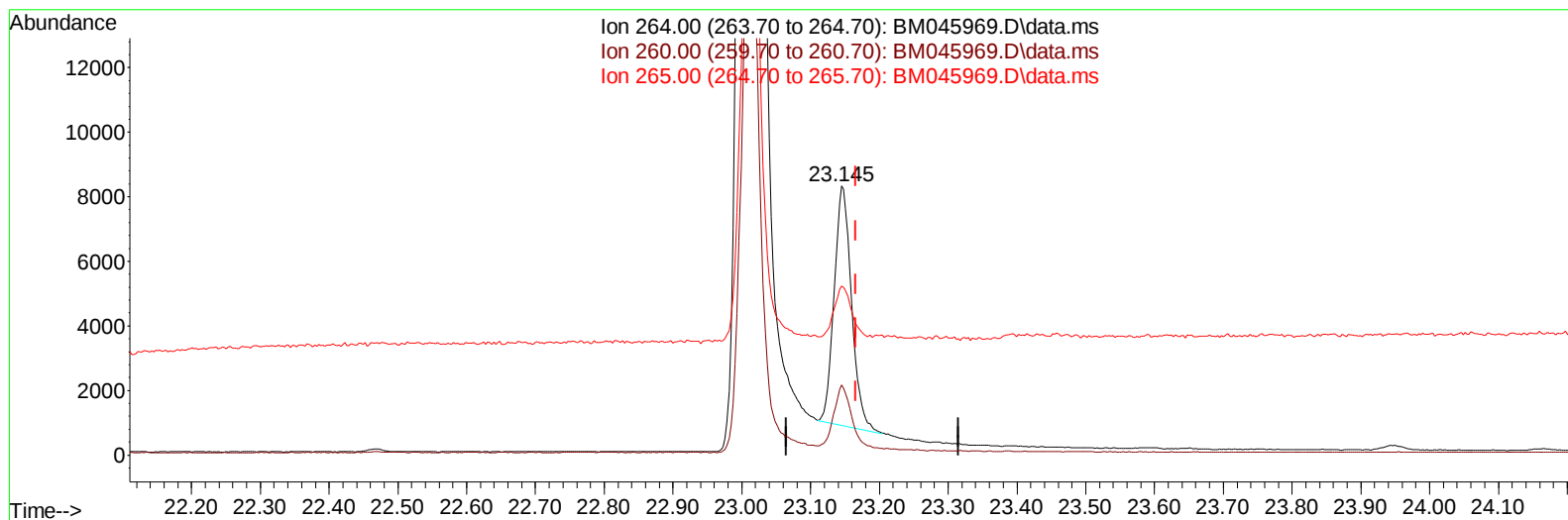
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061024\
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TIC: BM045969.D\data.ms

(23) Perylene-d12 (I)

23.145min (-0.020) 0.40 ng/ul m

response 13020

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	25.70	26.13
265.00	76.60	62.92
0.00	0.00	0.00

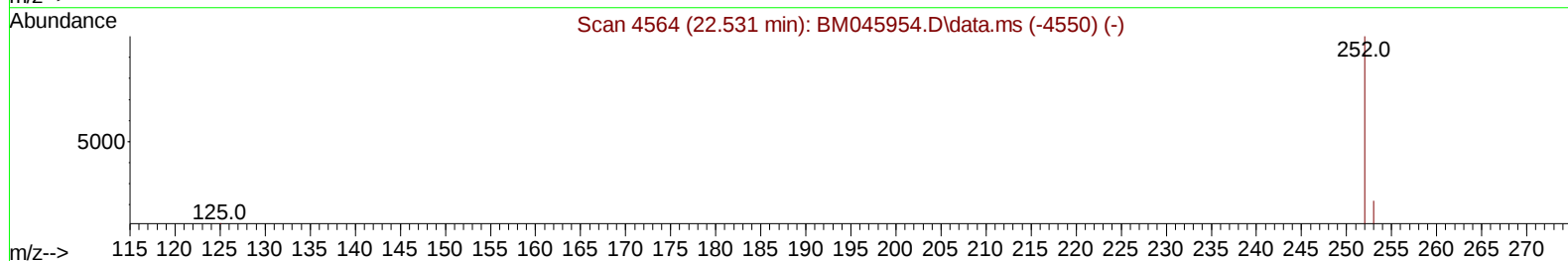
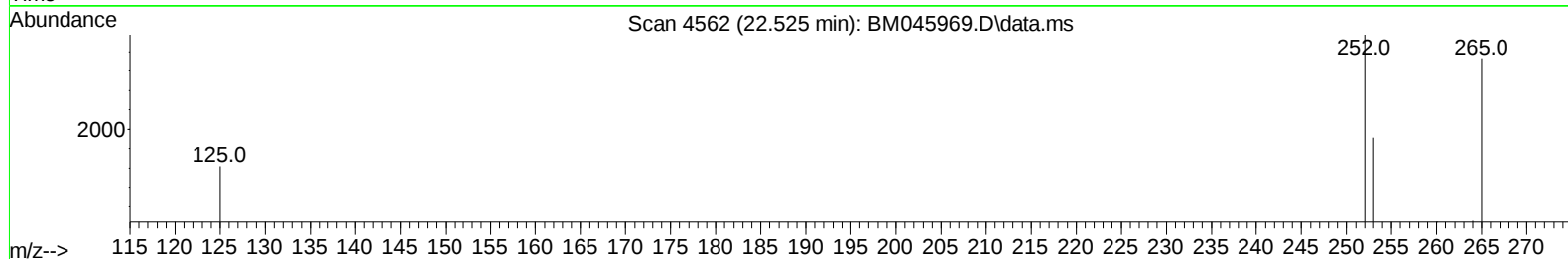
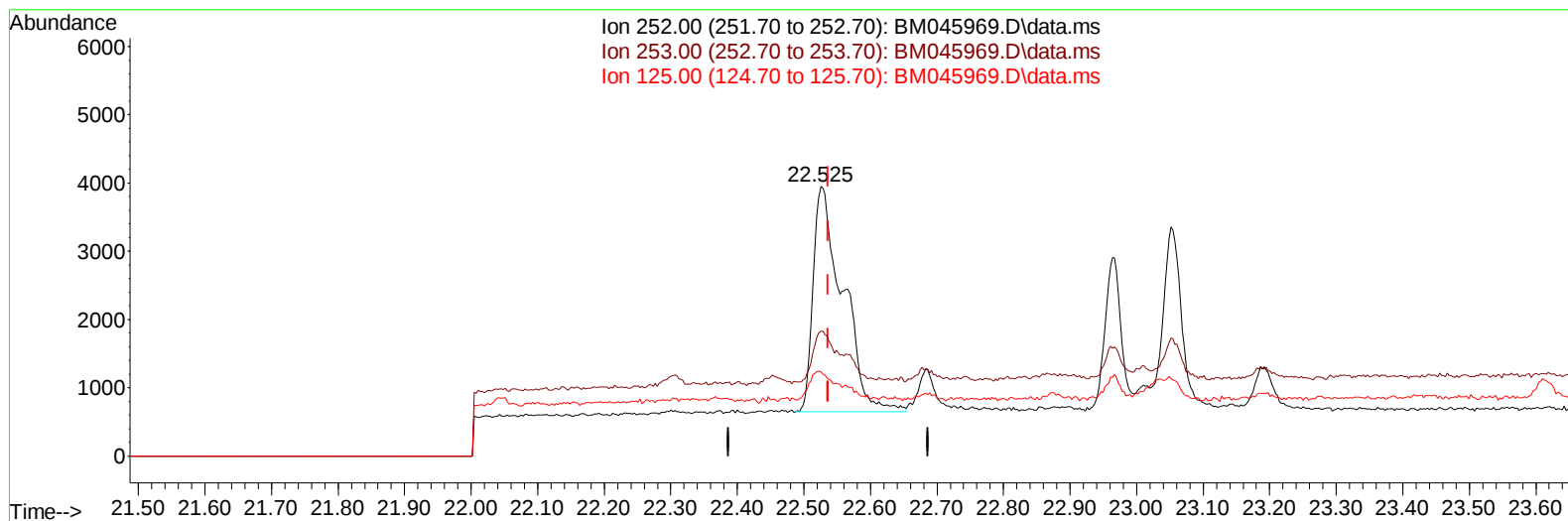
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Acq On : 11 Jun 2024 02:12
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Instrument :
BNA_M
ClientSampleId :
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TIC: BM045969.D\data.ms

(24) Benzo(b)fluoranthene

22.525min (-0.012) 0.23 ng/u1

response 10107

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.80	46.24
125.00	17.90	31.17
0.00	0.00	0.00

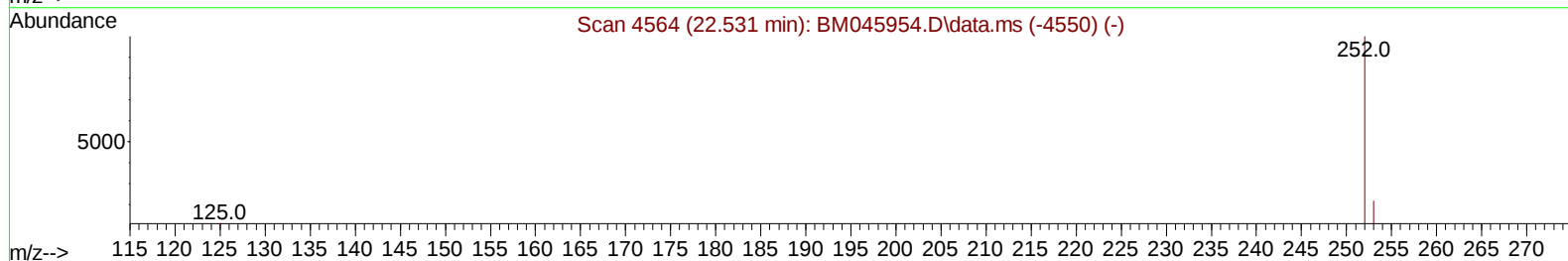
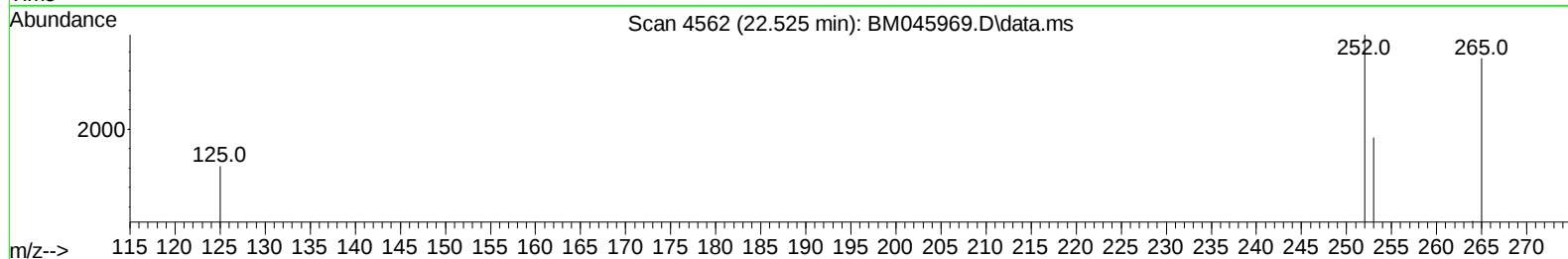
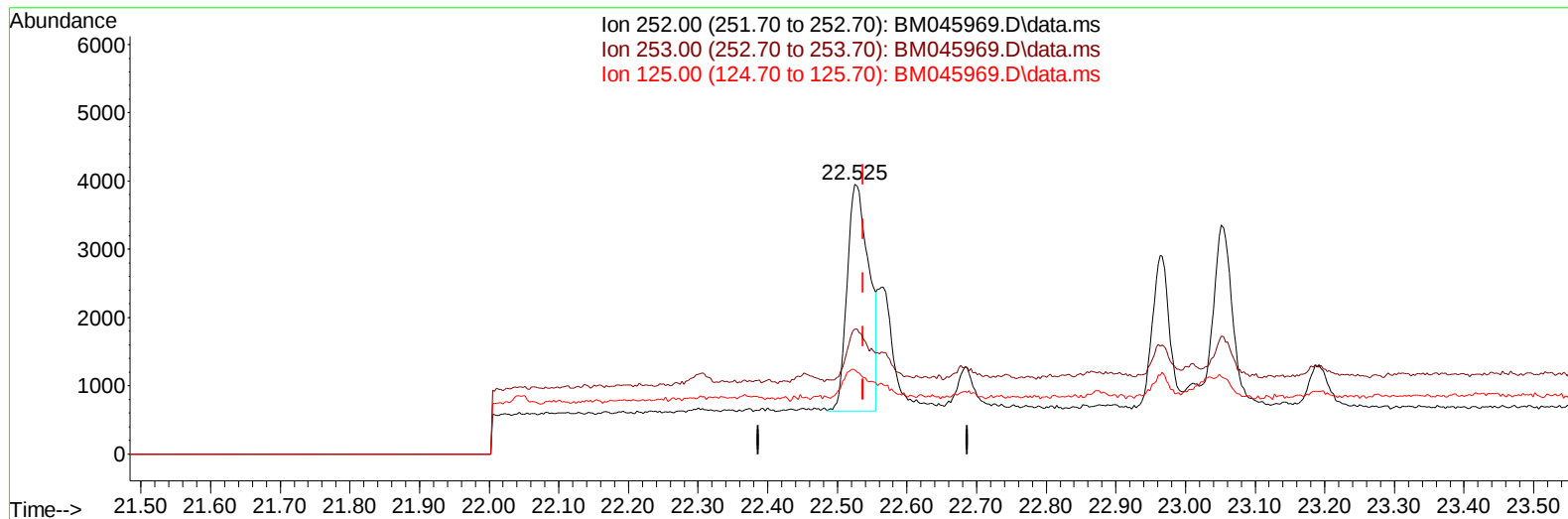
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061024\
Data File : BM045969.D
Acq On : 11 Jun 2024 02:12
Operator : MA/JU
Sample : P2642-17DL 5X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4B39DL

Manual IntegrationsAPPROVED

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Quant Time: Jun 11 02:42:13 2024
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TIC: BM045969.D\data.ms

(24) Benzo(b)fluoranthene

22.525min (-0.012) 0.17 ng/u1 m

response 7193

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.80	46.24
125.00	17.90	31.17
0.00	0.00	0.00

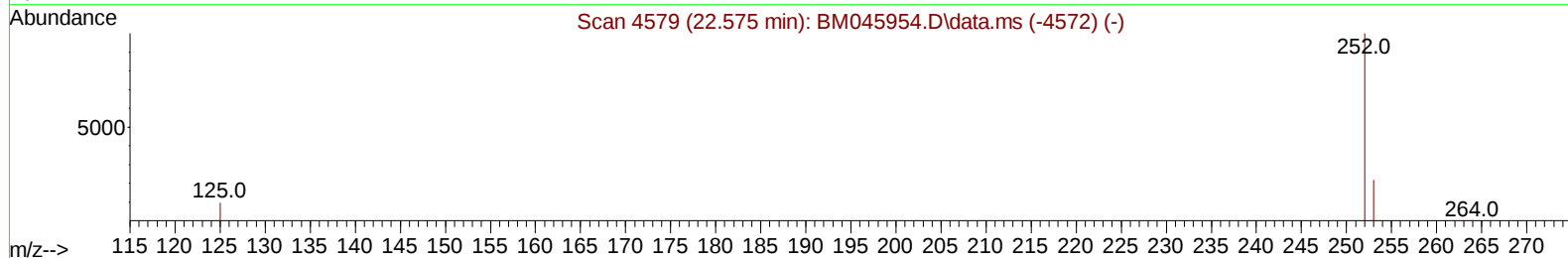
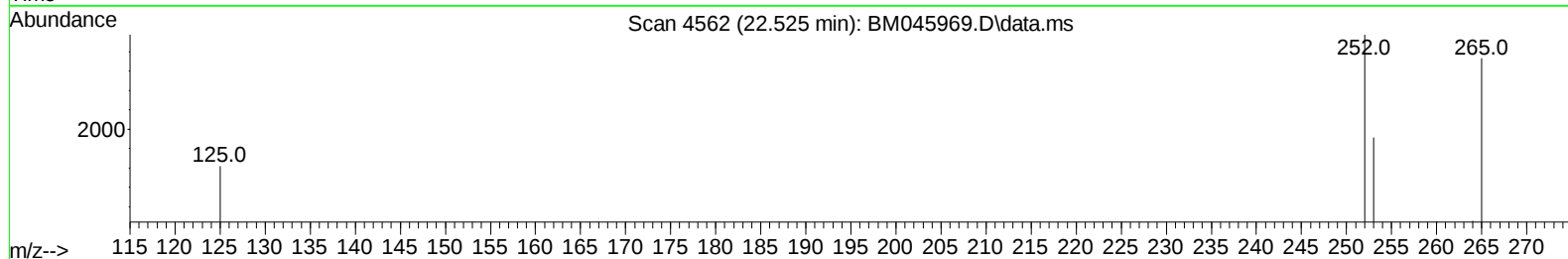
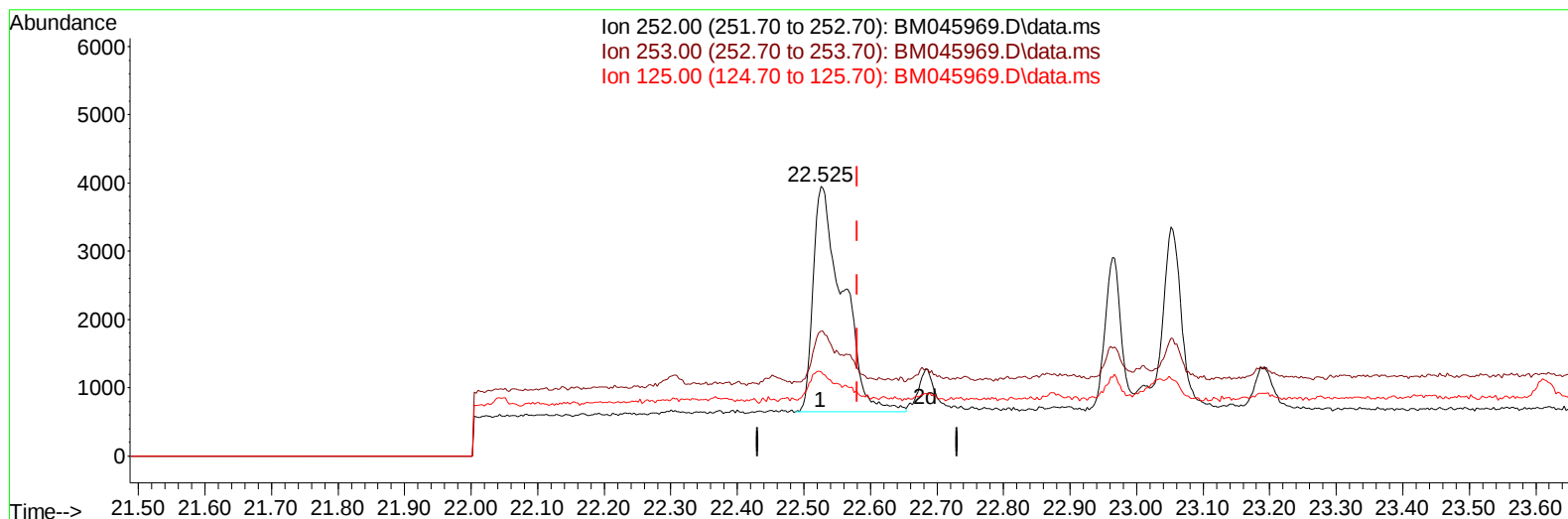
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061024\
Data File : BM045969.D
Acq On : 11 Jun 2024 02:12
Operator : MA/JU
Sample : P2642-17DL 5X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4B39DL

Manual IntegrationsAPPROVED

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TIC: BM045969.D\data.ms

(25) Benzo(k)fluoranthene

22.525min (-0.056) 0.21 ng/u1

response 10107

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.30	46.24#
125.00	17.70	31.17#
0.00	0.00	0.00

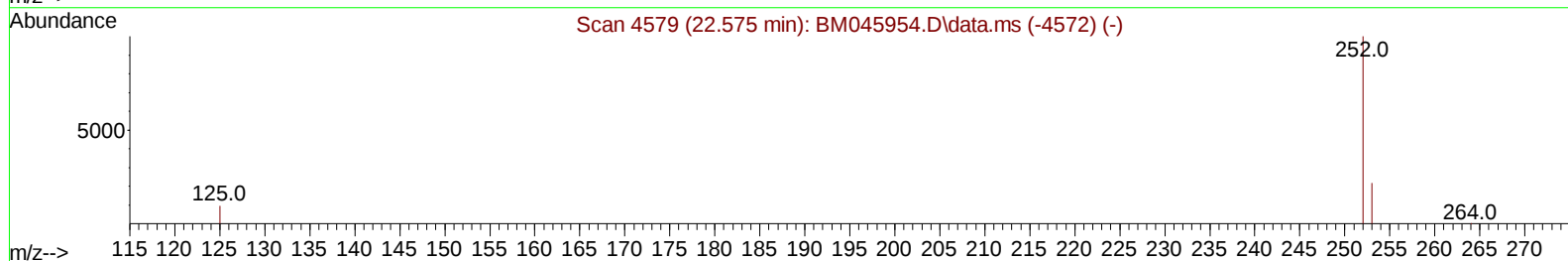
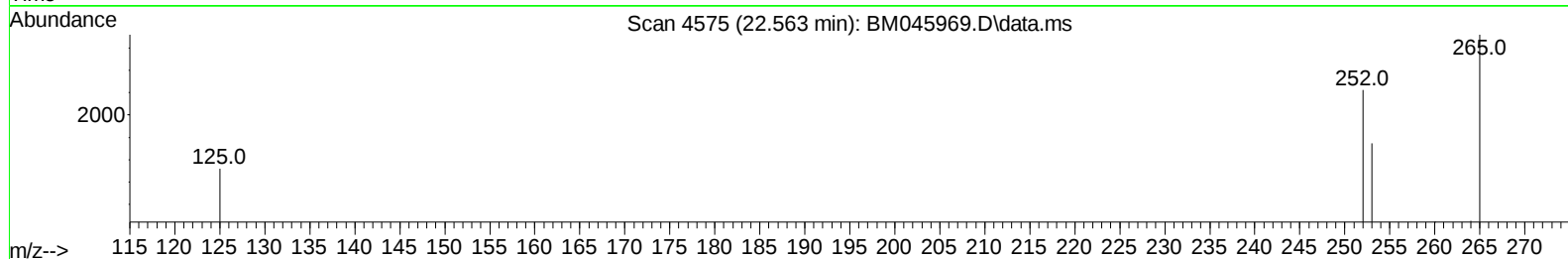
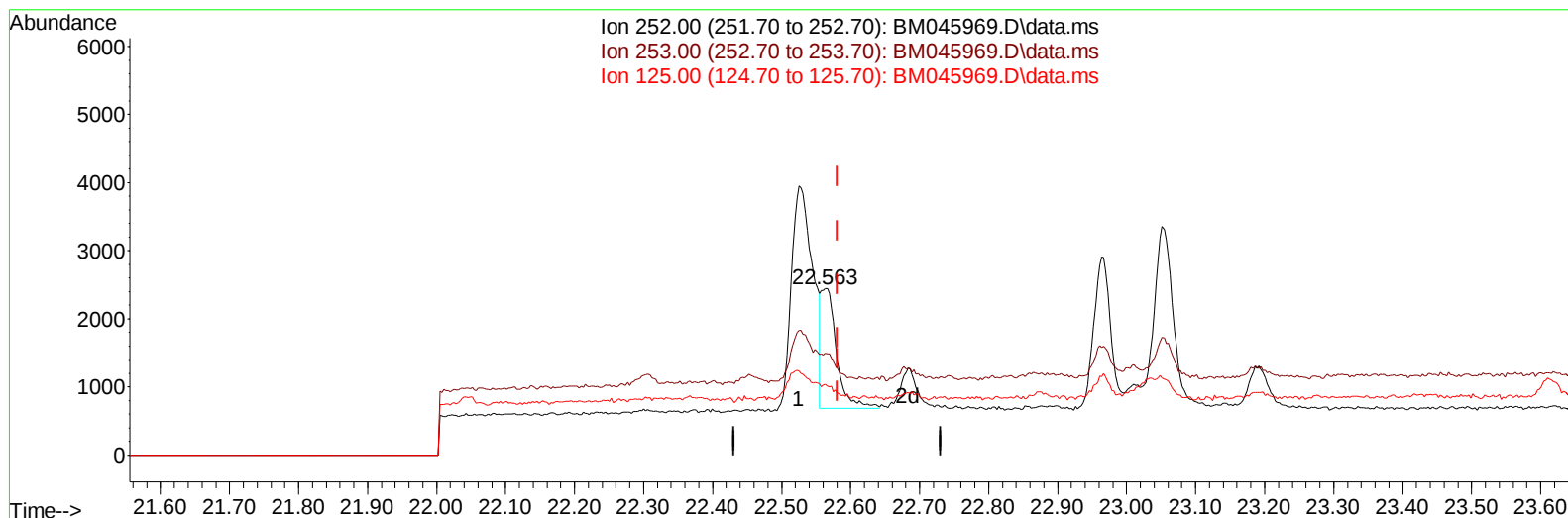
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061024\
 Data File : BM045969.D
 Acq On : 11 Jun 2024 02:12
 Operator : MA/JU
 Sample : P2642-17DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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TIC: BM045969.D\data.ms

(25) Benzo(k)fluoranthene

22.563min (-0.018) 0.06 ng/u1 m

response 2799

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.30	61.28#
125.00	17.70	42.49#
0.00	0.00	0.00

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 Operator : MA/JU
 Sample : P2642-17DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.459	152		4391	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.242	136		13538	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.108	164		7739	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.849	188		15184	0.400	ng/ul	-0.02
17) Chrysene-d12	21.044	240		11896	0.400	ng/ul	0.00
23) Perylene-d12	23.145	264		13020m	0.400	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.059	96		3823	0.764	ng/ul	-0.03
6) 2-Methylnaphthalene-d10	11.865	152		784	0.040	ng/ul	0.02
18) Fluoranthene-d10	18.880	212		1711	0.048	ng/ul	-0.01
Target Compounds							
							Qvalue
15) Phenanthrene	16.891	178		6543	0.153	ng/ul	97
16) Anthracene	16.984	178		1362	0.037	ng/ul#	81
19) Fluoranthene	18.913	202		11538	0.243	ng/ul	98
20) Pyrene	19.270	202		12071	0.244	ng/ul	98
21) Benzo(a)anthracene	21.029	228		4874	0.112	ng/ul#	92
22) Chrysene	21.082	228		5186	0.109	ng/ul	94
24) Benzo(b)fluoranthene	22.525	252		7193m	0.167	ng/ul	
25) Benzo(k)fluoranthene	22.563	252		2799m	0.059	ng/ul	
26) Benzo(a)pyrene	23.052	252		5059	0.118	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	25.216	276		3807	0.069	ng/ul#	93
29) Benzo(g,h,i)perylene	25.836	276		4115	0.086	ng/ul#	83

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

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