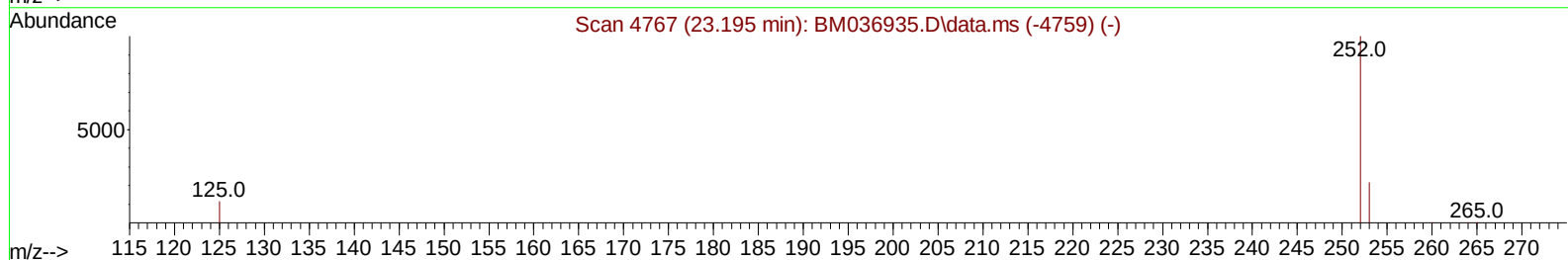
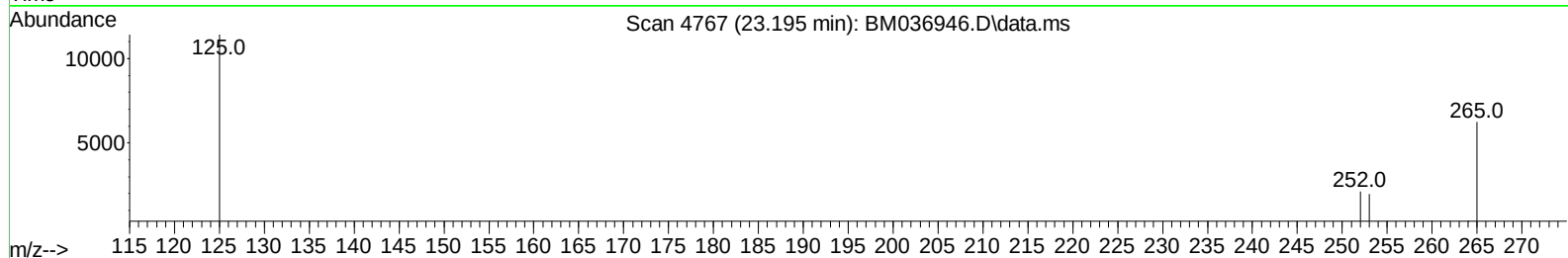
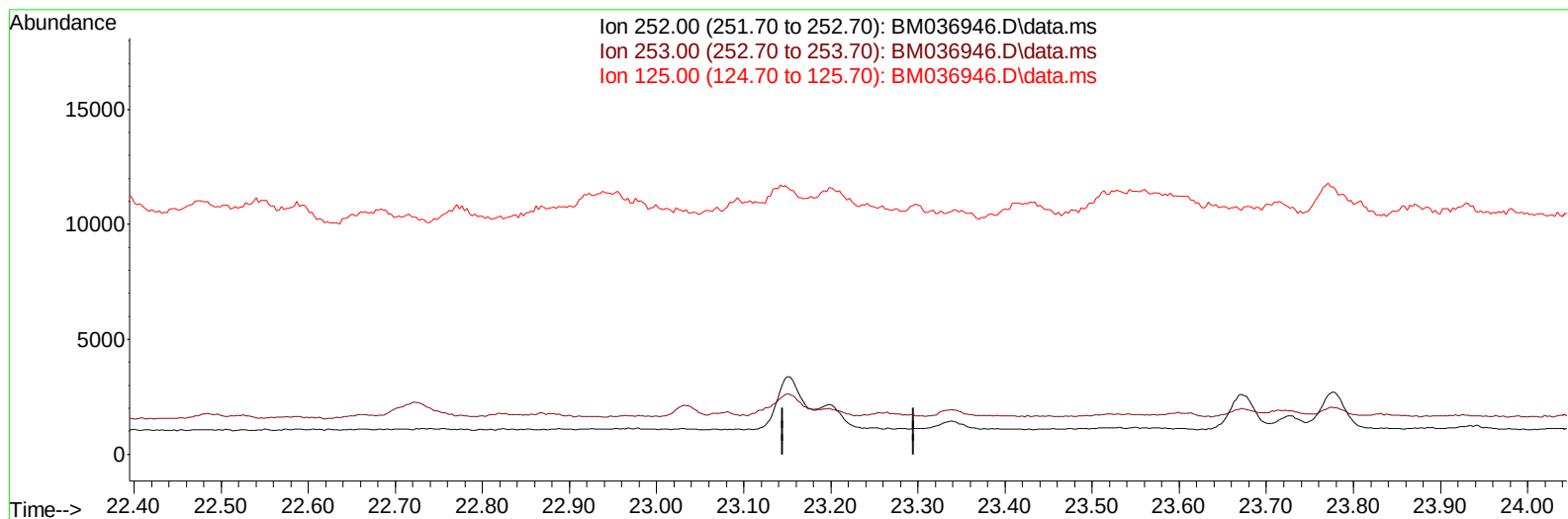


Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
Data File : BM036946.D
Acq On : 11 Oct 2022 15:16
Operator : CG/JU
Sample : N4824-04DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Quant Time: Oct 12 01:26:50 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 11 09:03:52 2022
Response via : Initial Calibration



TIC: BM036946.D\data.ms

(25) Benzo(k)fluoranthene

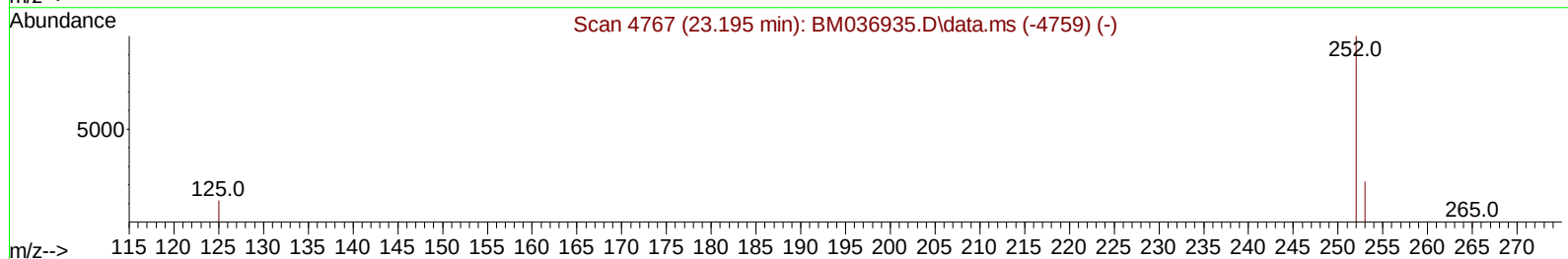
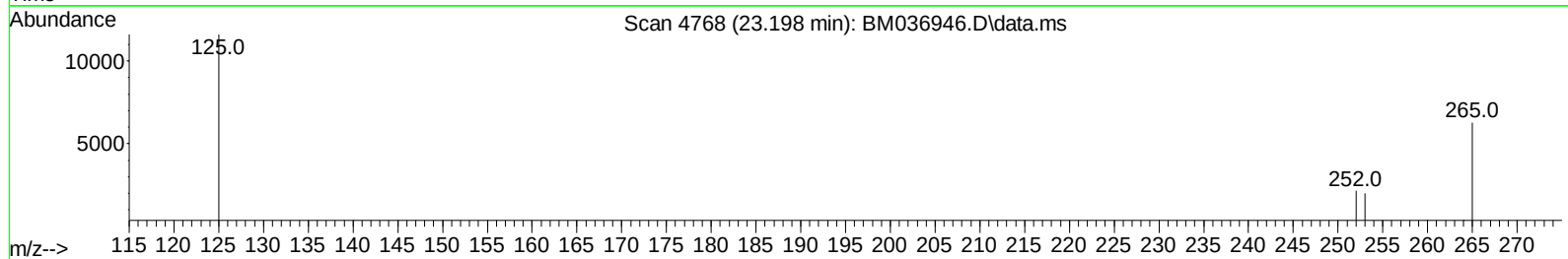
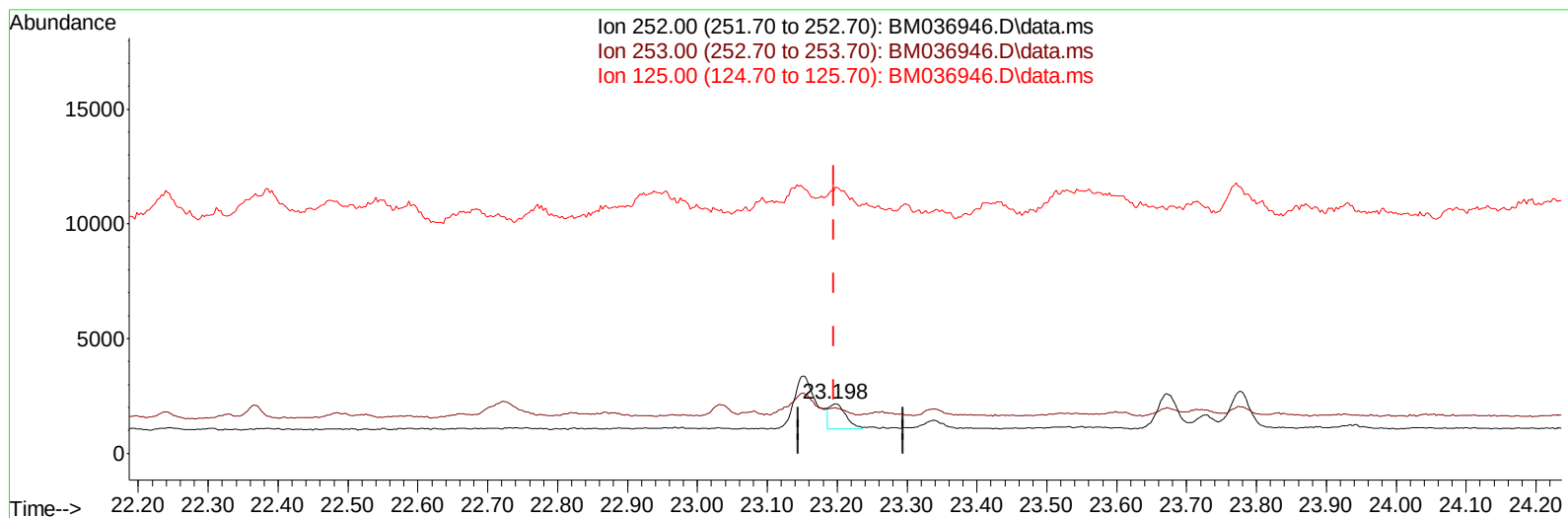
23.195min (-23.195) 0.00 ng/ul

response 0

Ion	Exp%	Act%
252.00	100.00	0.00
253.00	31.50	0.00#
125.00	22.90	0.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
Data File : BM036946.D
Acq On : 11 Oct 2022 15:16
Operator : CG/JU
Sample : N4824-04DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Quant Time: Oct 12 01:26:50 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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TIC: BM036946.D\data.ms

(25) Benzo(k)fluoranthene

23.198min (+ 0.003) 0.02 ng/ul m

response 1759

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.50	92.11#
125.00	22.90	536.06#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036946.D
 Acq On : 11 Oct 2022 15:16
 Operator : CG/JU
 Sample : N4824-04DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Quant Time: Oct 12 01:26:50 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 09:03:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.955	152	8497	0.400	ng/ul	0.00
4) Naphthalene-d8	10.759	136	27952	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.579	164	14555	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.317	188	27837	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.485	240	21649	0.400	ng/ul	# 0.00
23) Perylene-d12	23.885	264	21622	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	8977	0.749	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.343	152	2509	0.059	ng/ul	0.00
18) Fluoranthene-d10	19.336	212	4240	0.065	ng/ul	0.00
Target Compounds						
					Qvalue	
8) 1-Methylnaphthalene	12.635	142	1035	0.021	ng/ul	99
10) Acenaphthylene	14.302	152	2271	0.038	ng/ul#	18
11) Acenaphthene	14.640	153	49391	0.941	ng/ul	99
12) Fluorene	15.625	166	20182	0.339	ng/ul	99
15) Phenanthrene	17.360	178	6058	0.068	ng/ul#	77
16) Anthracene	17.448	178	14862	0.200	ng/ul#	93
19) Fluoranthene	19.368	202	79231	0.920	ng/ul	98
20) Pyrene	19.726	202	62032	0.691	ng/ul	98
21) Benzo(a)anthracene	21.470	228	10266	0.131	ng/ul#	87
22) Chrysene	21.520	228	8572	0.102	ng/ul#	84
24) Benzo(b)fluoranthene	23.151	252	5169	0.054	ng/ul#	1
26) Benzo(a)pyrene	23.777	252	3050	0.039	ng/ul#	1

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

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