

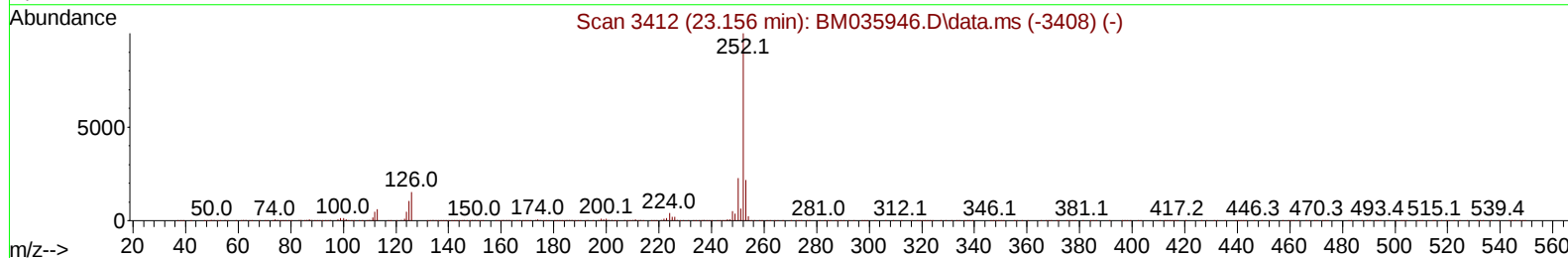
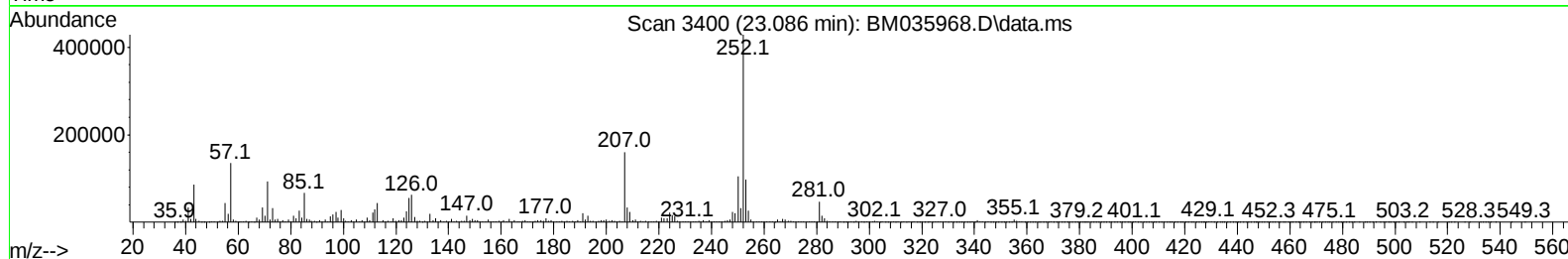
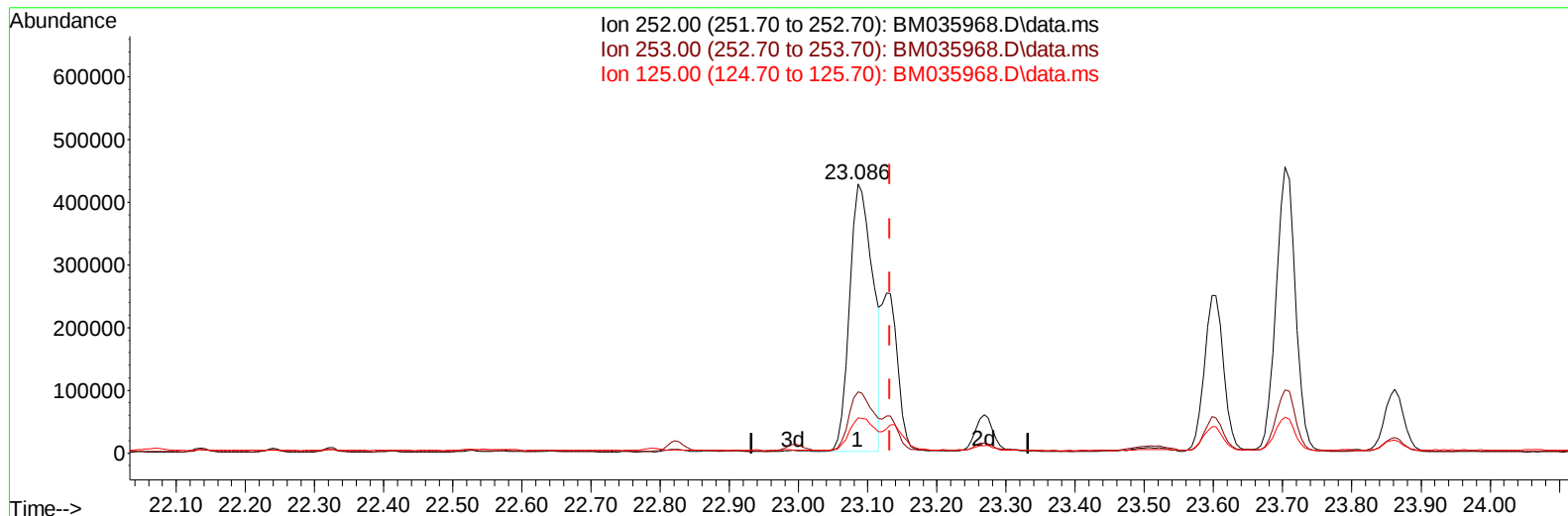
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
Data File : BM035968.D
Acq On : 28 Jul 2022 02:19
Operator : CG/JU
Sample : N3741-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
H0B36

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/28/2022
Supervised By :mohammad ahmed 07/29/2022

Quant Time: Jul 28 03:13:50 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 27 23:55:50 2022
Response via : Initial Calibration



TIC: BM035968.D\data.ms

(91) Benzo(k)fluoranthene

23.086min (-0.047) 11.66 ng/u1

response 999470

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.70	22.73
125.00	13.20	13.04
0.00	0.00	0.00

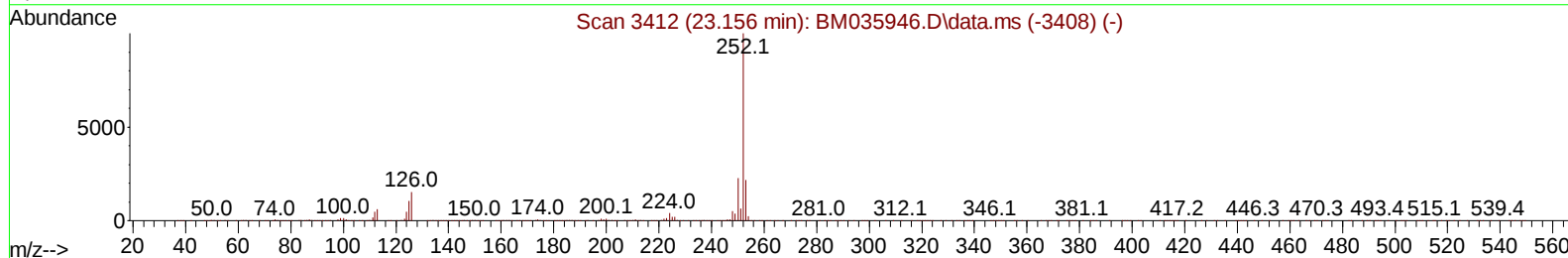
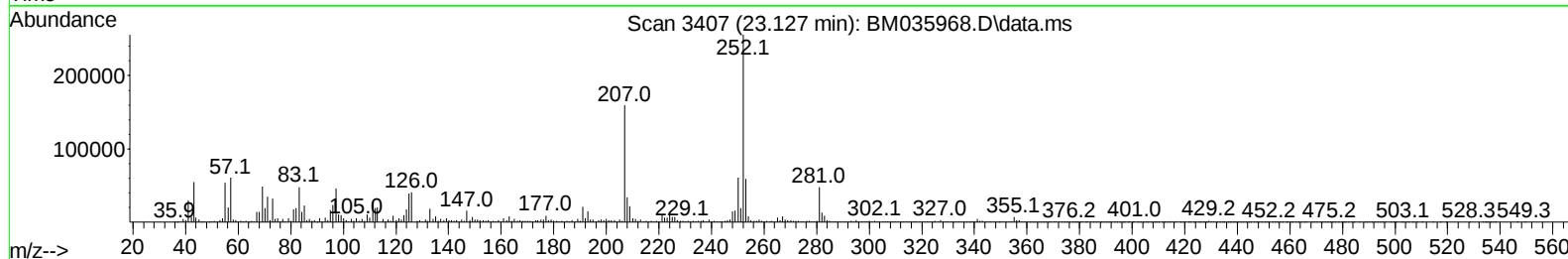
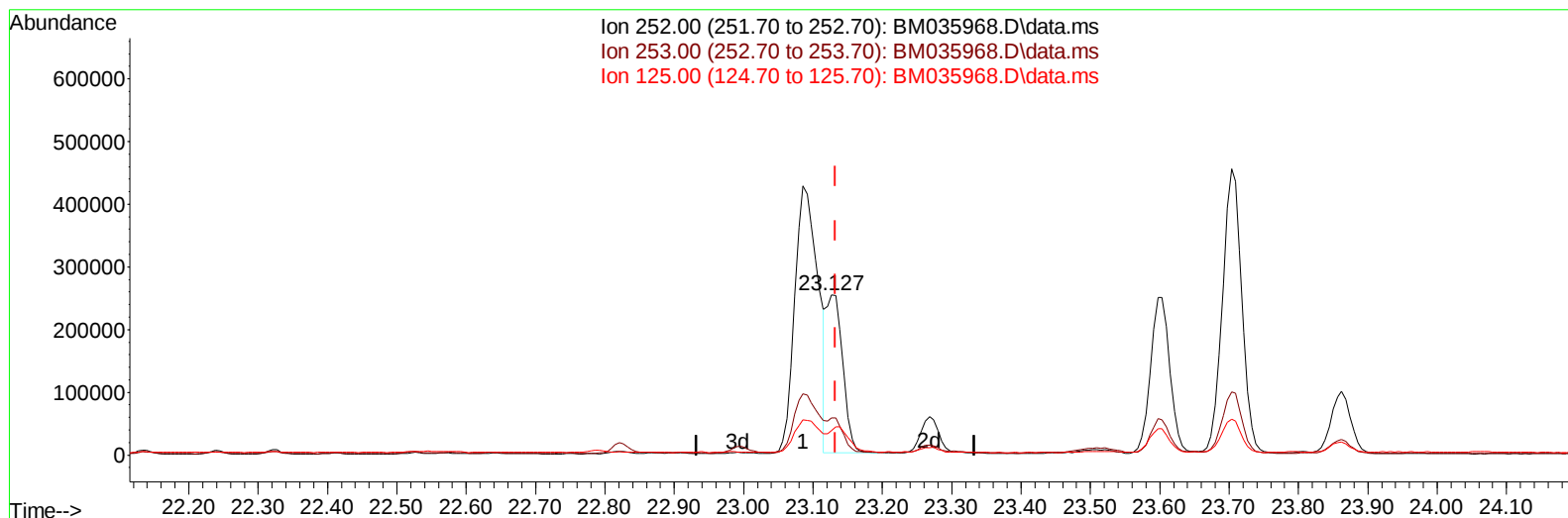
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035968.D
 Acq On : 28 Jul 2022 02:19
 Operator : CG/JU
 Sample : N3741-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0B36

Manual IntegrationsAPPROVED

Quant Time: Jul 28 03:13:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 23:55:50 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/28/2022
 Supervised By :mohammad ahmed 07/29/2022



TIC: BM035968.D\data.ms

(91) Benzo(k)fluoranthene

23.127min (-0.006) 4.75 ng/ul m

response 407118

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.70	23.04
125.00	13.20	15.16
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035968.D
 Acq On : 28 Jul 2022 02:19
 Operator : CG/JU
 Sample : N3741-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0B36

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/28/2022
 Supervised By :mohammad ahmed 07/29/2022

Quant Time: Jul 28 03:13:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 23:55:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	363232	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1555170	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	884880	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1632242	20.000	ng/ul	0.00
79) Chrysene-d12	21.439	240	1316930	20.000	ng/ul	0.00
88) Perylene-d12	23.809	264	1353025	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	28243	2.926	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.051	99	568419	18.661	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	372495	18.344	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	502033	19.856	ng/ul	0.00
15) 4-Methylphenol-d8	8.598	113	361289	14.641	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	256472	20.812	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	266220	22.173	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	438369	20.188	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	447524	12.174	ng/ul	0.00
46) Dimethylphthalate-d6	13.933	166	1330788	22.078	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1714671	22.296	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	158138	12.776	ng/ul	0.00
60) Fluorene-d10	15.510	176	1160003	21.297	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	114035	12.997	ng/ul	0.00
73) Anthracene-d10	17.357	188	1575424	21.349	ng/ul	0.00
81) Pyrene-d10	19.651	212	1723263	23.611	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.656	264	1528649	22.254	ng/ul	0.00
Target Compounds						
					Qvalue	
52) Acenaphthene	14.580	153	65725	1.191	ng/ul	99
72) Phenanthrene	17.304	178	1140755	13.040	ng/ul	99
74) Anthracene	17.392	178	180987	2.062	ng/ul	98
77) Carbazole	17.663	167	142920	1.763	ng/ul	99
80) Fluoranthene	19.315	202	1657519	19.008	ng/ul	97
82) Pyrene	19.680	202	1850313	20.586	ng/ul	95
85) Benzo(a)anthracene	21.421	228	771872	9.312	ng/ul	99
87) Chrysene	21.474	228	996892	12.326	ng/ul	99
90) Benzo(b)fluoranthene	23.086	252	999470	11.069	ng/ul	99
91) Benzo(k)fluoranthene	23.127	252	407118m	4.751	ng/ul	
93) Benzo(a)pyrene	23.703	252	879285	10.223	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	26.274	276	574837	6.270	ng/ul#	91
95) Dibenzo(a,h)anthracene	26.285	278	158629	2.017	ng/ul	94
96) Benzo(g,h,i)perylene	27.038	276	467091	6.148	ng/ul	95

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035968.D
 Acq On : 28 Jul 2022 02:19
 Operator : CG/JU
 Sample : N3741-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0B36

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/28/2022
 Supervised By :mohammad ahmed 07/29/2022

Quant Time: Jul 28 03:13:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 23:55:50 2022
 Response via : Initial Calibration

