

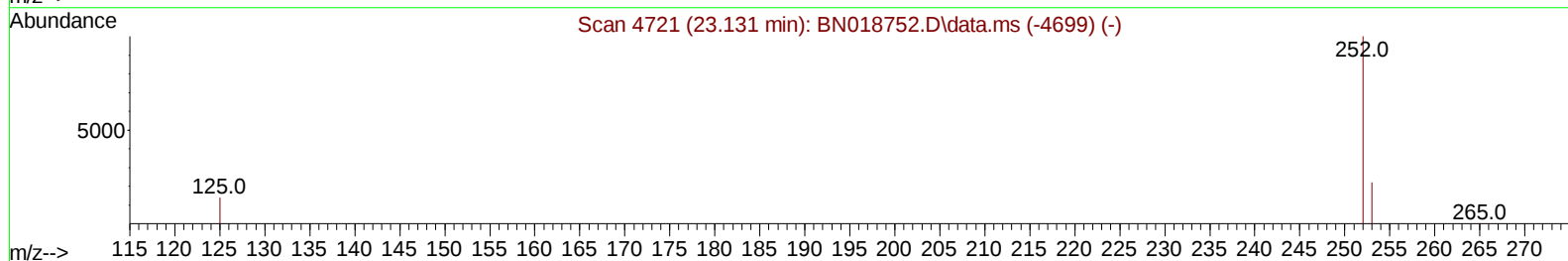
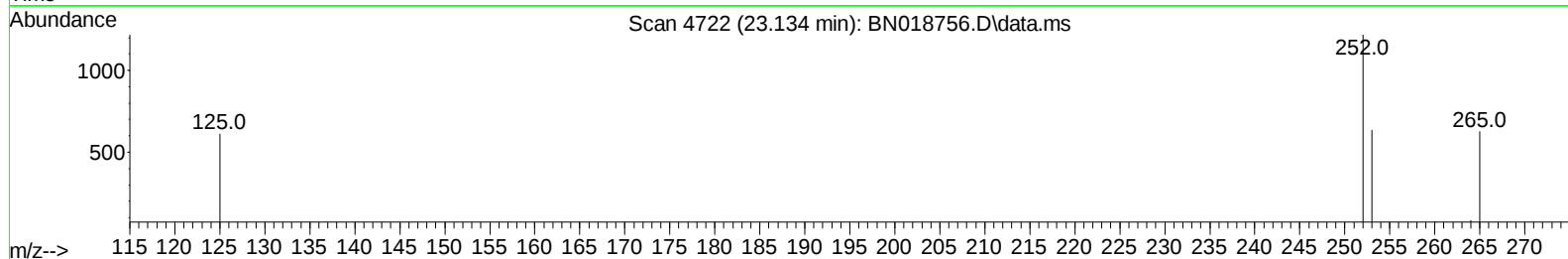
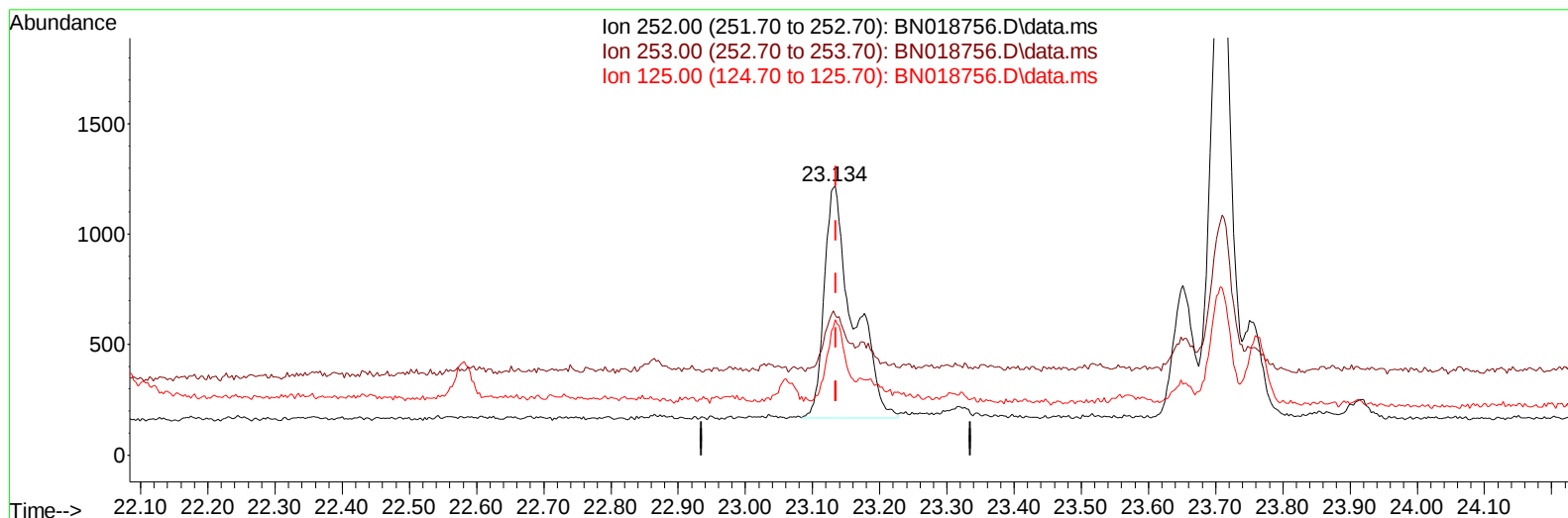
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022822\
 Data File : BN018756.D
 Acq On : 28 Feb 2022 12:40
 Operator : CG/JU
 Sample : N1656-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGQ11

Manual IntegrationsAPPROVED

Quant Time: Feb 28 23:54:50 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN021822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 18 13:39:55 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/01/2022
 Supervised By :Yogesh Patel 03/02/2022



TIC: BN018756.D\data.ms

(24) Benzo(b)fluoranthene

23.134min (-0.002) 0.06 ng/u1

response 3154

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	23.20	52.05#
125.00	8.90	50.25#
0.00	0.00	0.00

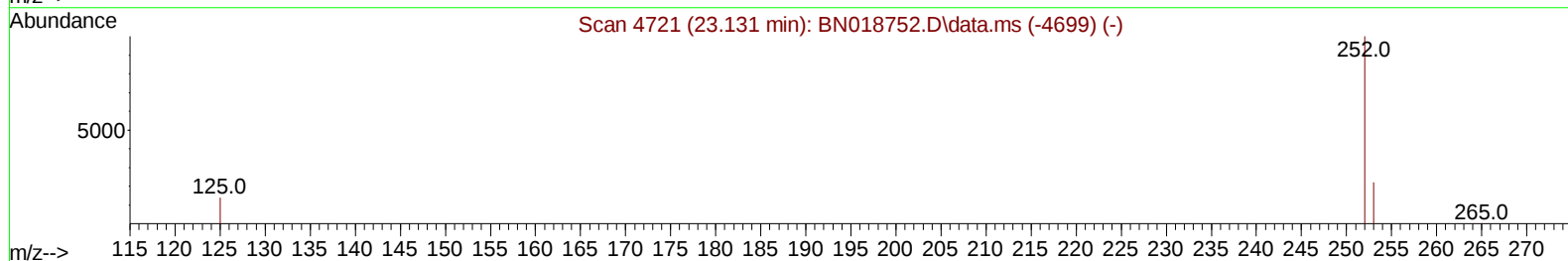
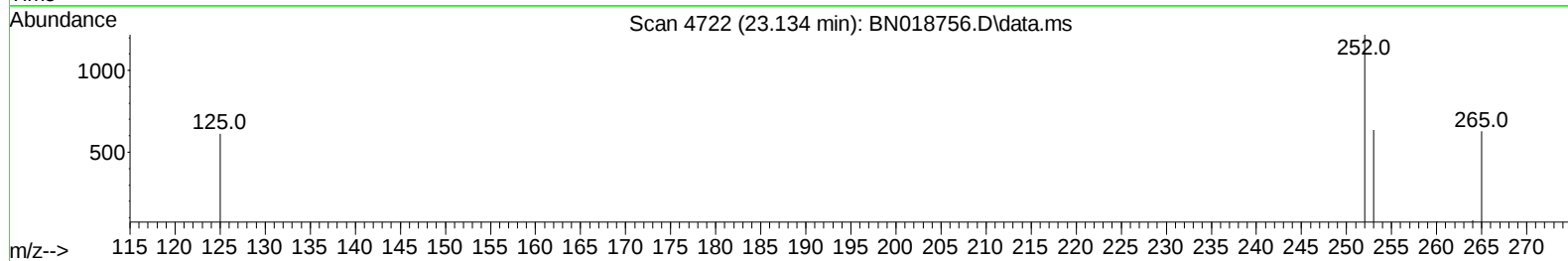
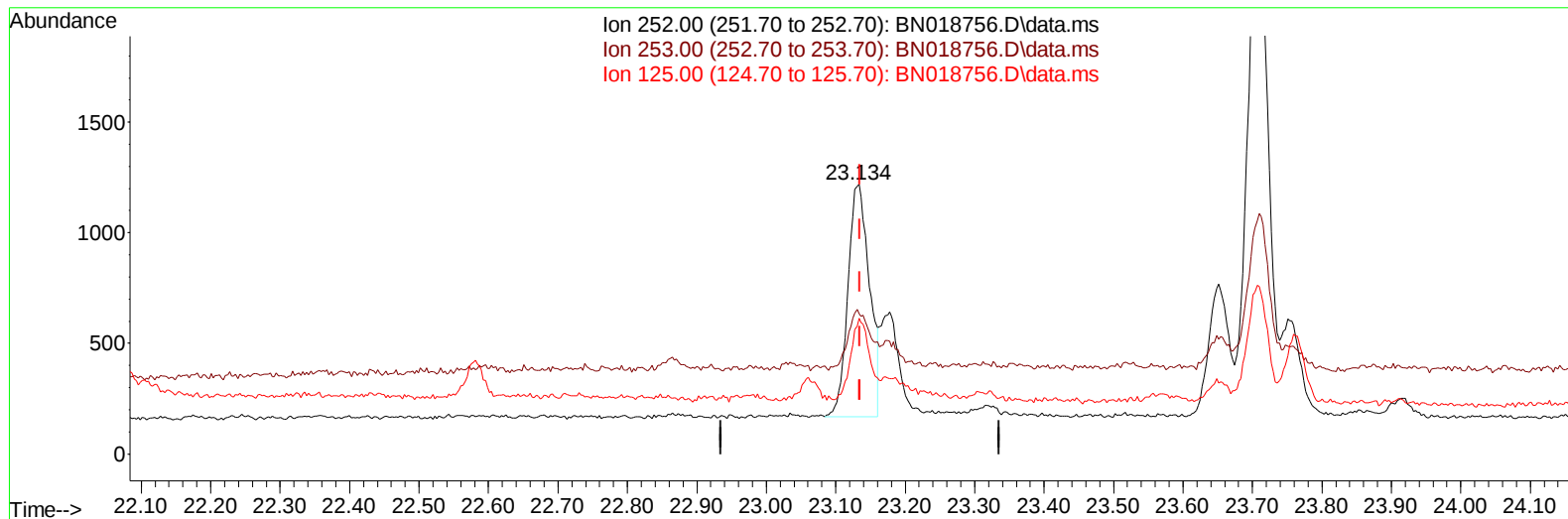
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TIC: BN018756.D\data.ms

(24) Benzo(b)fluoranthene

23.134min (-0.002) 0.05 ng/u1 m

response 2264

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	23.20	52.05#
125.00	8.90	50.25#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.913	152	5845	0.400	ng/ul	0.00
4) Naphthalene-d8	10.721	136	13966	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.547	164	7082	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.284	188	13594	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.465	240	11308	0.400	ng/ul	# 0.00
23) Perylene-d12	23.861	264	11091	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	36571	5.509	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.310	152	6802	0.356	ng/ul	0.00
18) Fluoranthene-d10	19.312	212	16132	0.478	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.322	88	661	0.097	ng/ul#	83
15) Phenanthrene	17.326	178	1910	0.043	ng/ul#	87
19) Fluoranthene	19.340	202	2575	0.054	ng/ul#	81
20) Pyrene	19.703	202	1847	0.038	ng/ul#	71
22) Chrysene	21.500	228	1463	0.032	ng/ul#	84
24) Benzo(b)fluoranthene	23.134	252	2264m	0.046	ng/ul	
27) Indeno(1,2,3-cd)pyrene	26.343	276	1212	0.022	ng/ul#	78
29) Benzo(g,h,i)perylene	27.104	276	1375	0.029	ng/ul#	66

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

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