

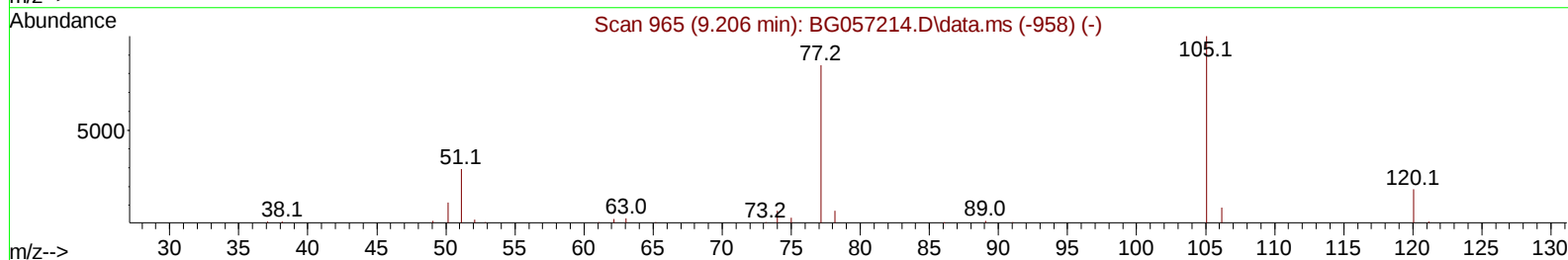
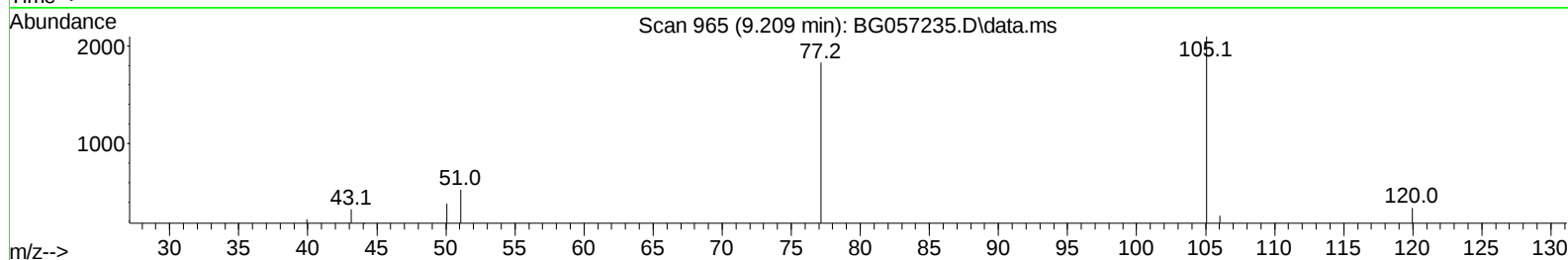
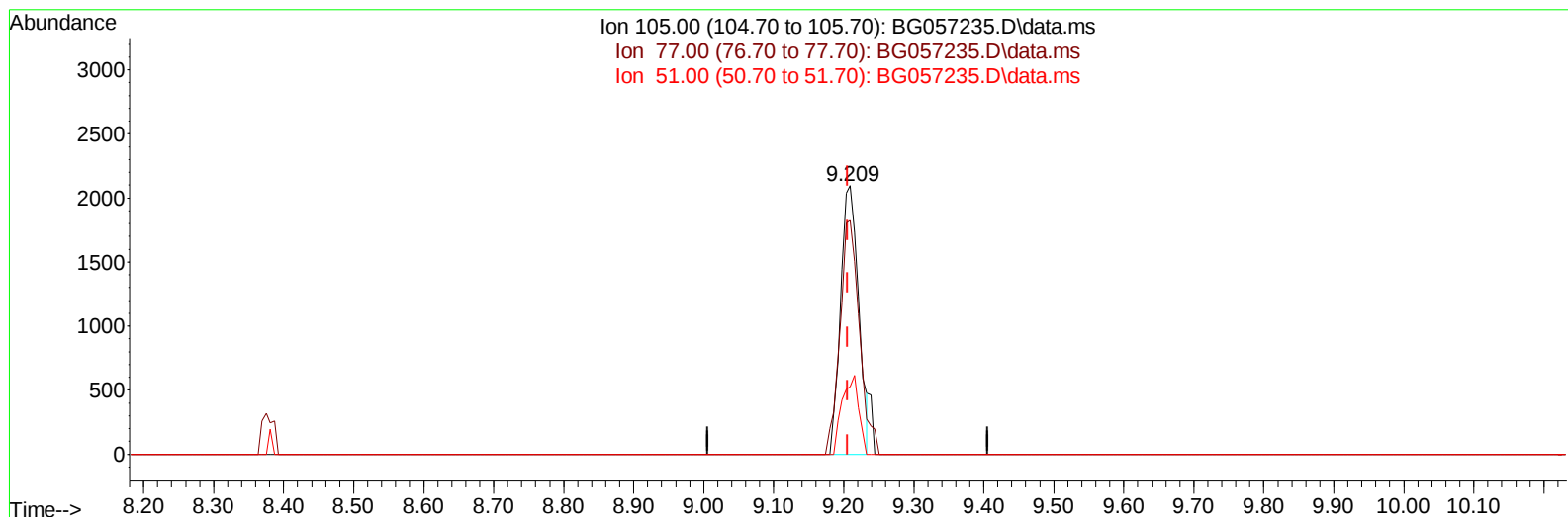
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
 Data File : BG057235.D
 Acq On : 29 Apr 2023 3:04
 Operator : CG/JU
 Sample : 02417-09
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW8X5

Manual IntegrationsAPPROVED

Quant Time: Apr 29 04:21:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 28 23:37:02 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/01/2023
 Supervised By :Jagrut Upadhyay 05/01/2023



TIC: BG057235.D\data.ms

(16) Acetophenone

9.209min (+ 0.003) 2.46 ng/u1

response 3755

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	88.20	87.21
51.00	30.60	25.01
0.00	0.00	0.00

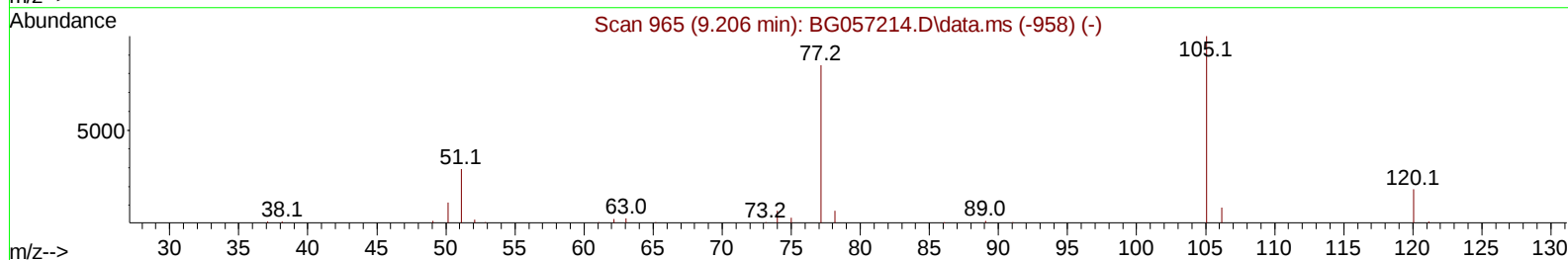
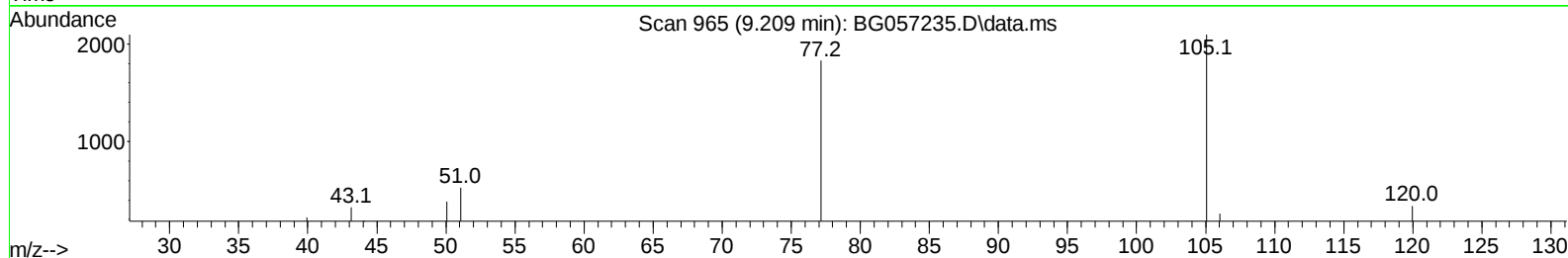
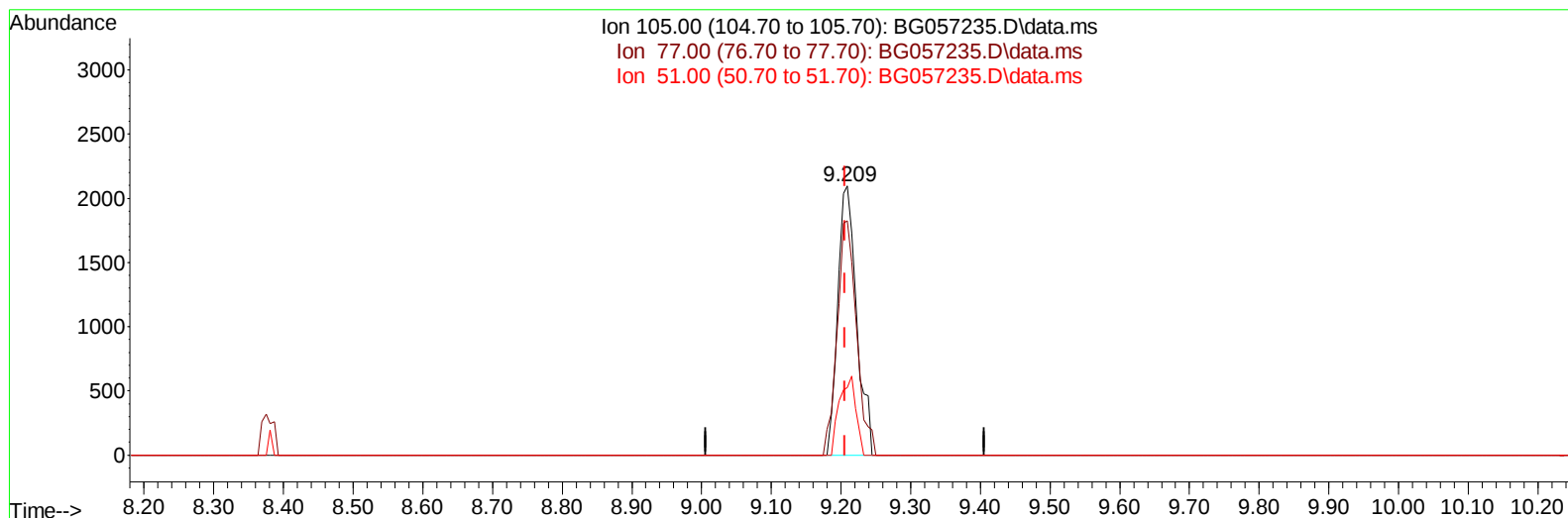
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TIC: BG057235.D\data.ms

(16) Acetophenone

9.209min (+ 0.003) 2.56 ng/ul m

response 3918

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	88.20	87.21
51.00	30.60	25.01
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.381	152		12365	20.000	ng/ul	0.00
20) Naphthalene-d8	11.212	136		52378	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.990	164		44521	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.727	188		115116	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.033	240		102339	20.000	ng/ul	0.00
88) Perylene-d12	25.540	264		106571	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.652	96		414	1.473	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.512	99		14833	12.685	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.682	67		8959	13.026	ng/ul	0.00
11) 2-Chlorophenol-d4	7.905	132		10748	13.947	ng/ul	0.00
15) 4-Methylphenol-d8	9.074	113		9559	10.755	ng/ul	0.00
21) Nitrobenzene-d5	9.550	128		5444	13.051	ng/ul	0.00
24) 2-Nitrophenol-d4	10.284	143		6674	13.710	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.831	165		12970	13.810	ng/ul	0.00
31) 4-Chloroaniline-d4	11.342	131		14187	11.397	ng/ul	0.00
46) Dimethylphthalate-d6	14.373	166		49357	13.991	ng/ul	0.00
49) Acenaphthylene-d8	14.684	160		54675	13.790	ng/ul	0.00
54) 4-Nitrophenol-d4	15.178	143		5046	9.827	ng/ul	0.00
60) Fluorene-d10	15.971	176		46674	14.203	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.088	200		6401	9.385	ng/ul	0.00
73) Anthracene-d10	17.827	188		77922	14.830	ng/ul	0.00
81) Pyrene-d10	20.089	212		94525	15.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.293	264		83976	15.467	ng/ul	0.00
Target Compounds							
16) Acetophenone	9.209	105		3918m	2.562	ng/ul	Qvalue

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

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