

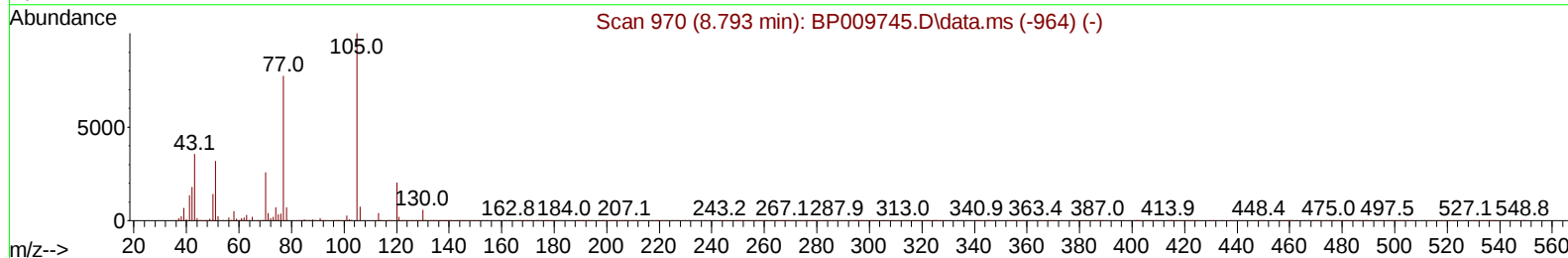
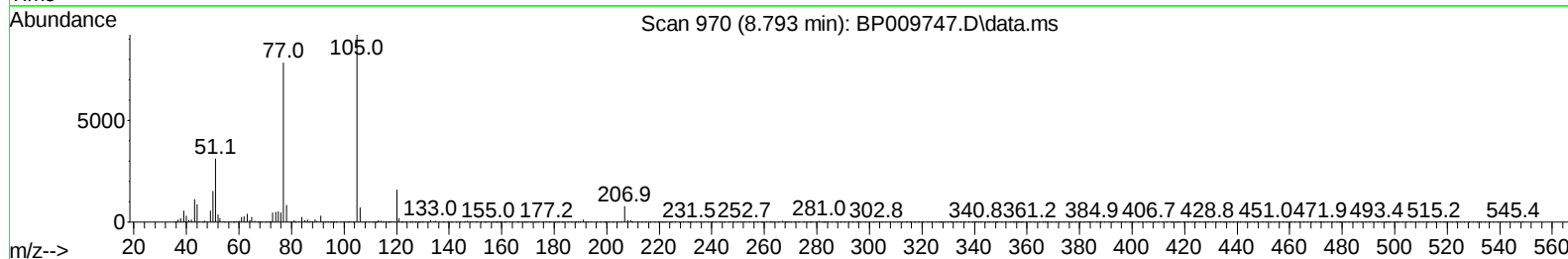
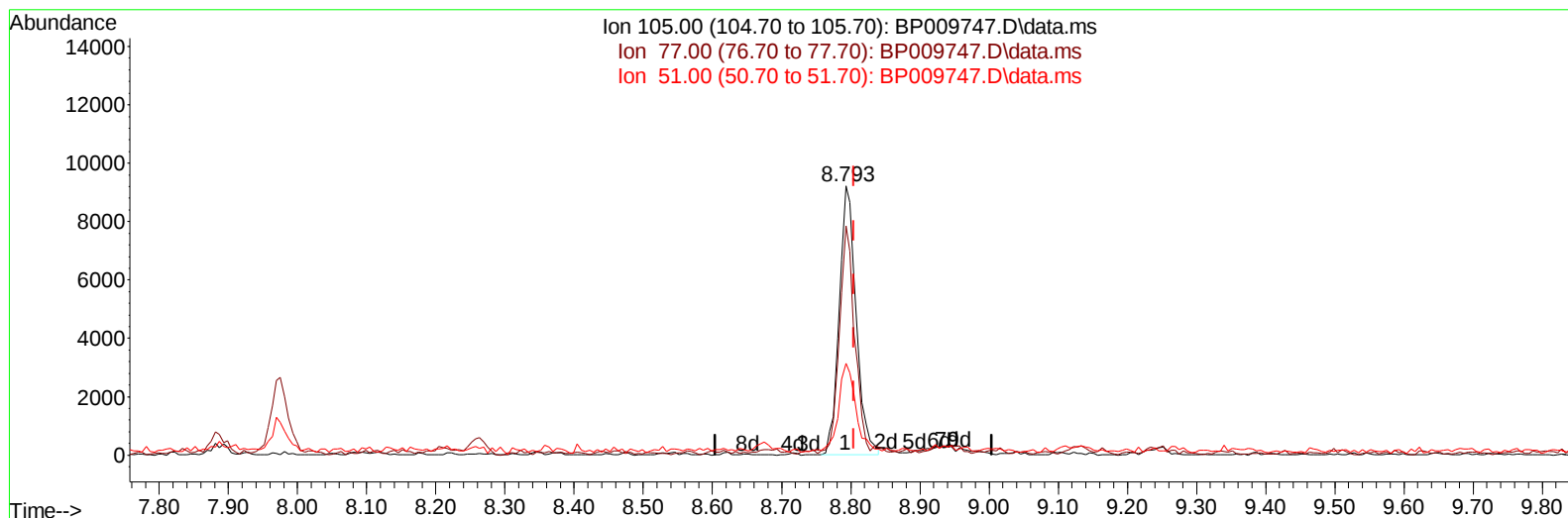
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009747.D
 Acq On : 11 Apr 2022 12:23
 Operator : CG/JU
 Sample : N2182-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R14

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:06:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/12/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009747.D\data.ms

(16) Acetophenone

8.793min (-0.012) 1.29 ng/u1

response 15687

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	85.15#
51.00	17.30	34.06#
0.00	0.00	0.00

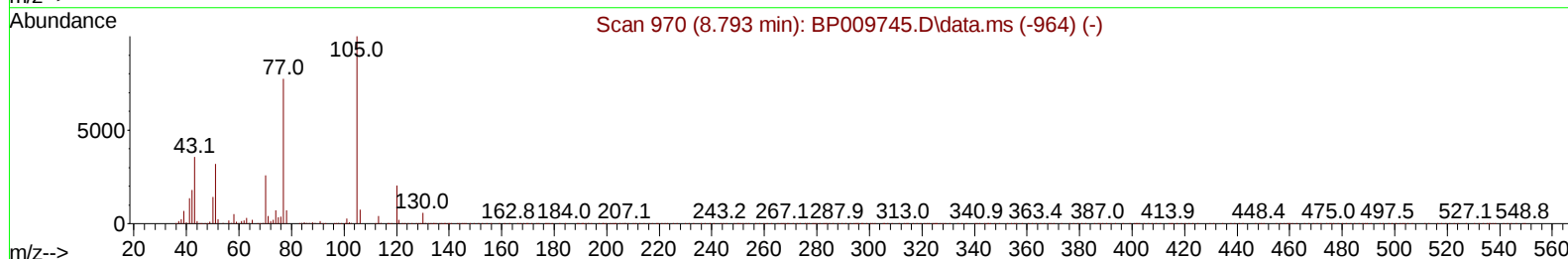
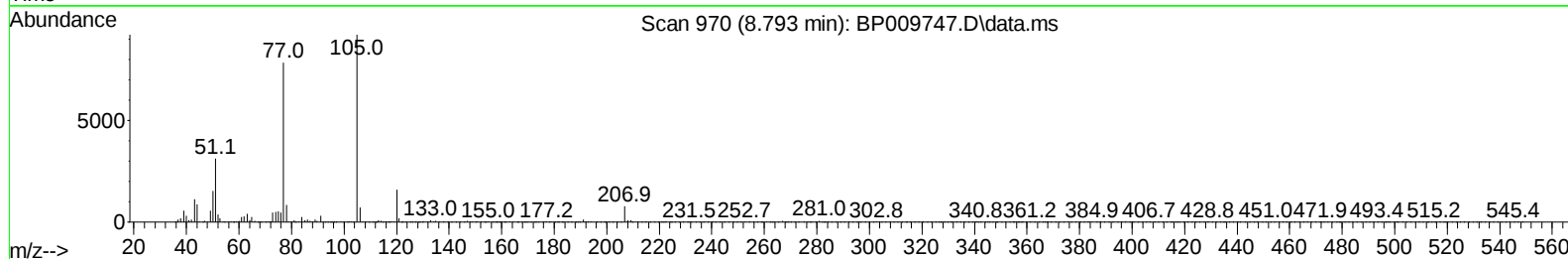
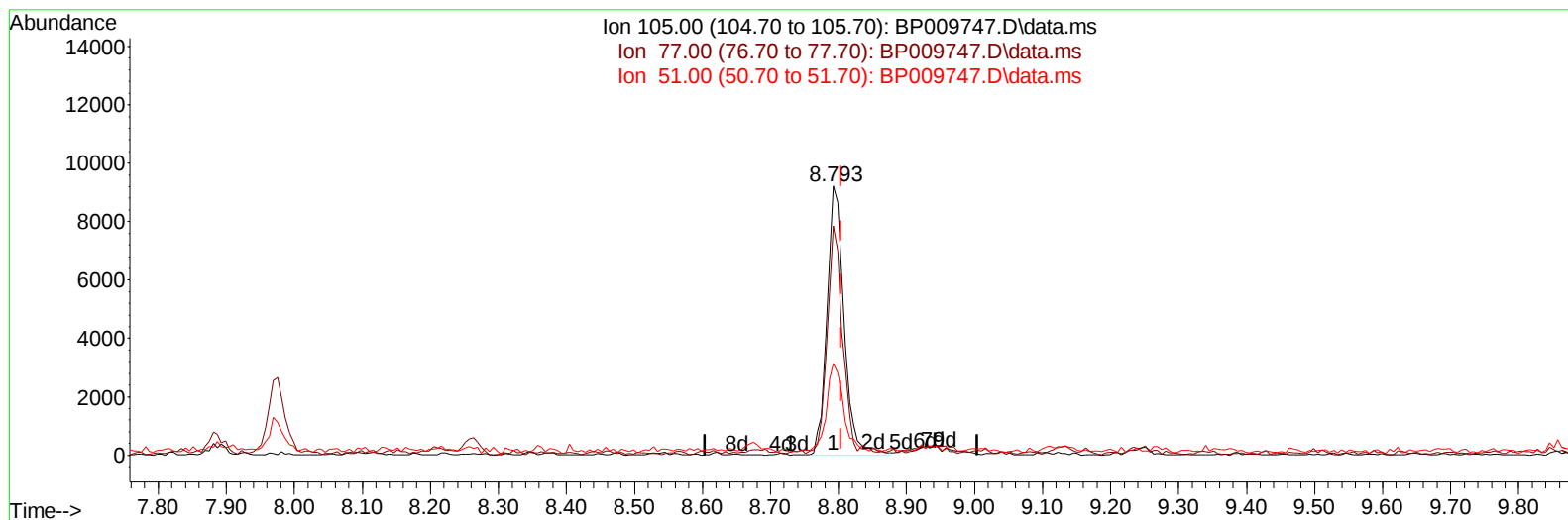
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TIC: BP009747.D\data.ms

(16) Acetophenone

8.793min (-0.012) 1.32 ng/ul m

response 16065

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	85.15#
51.00	17.30	34.06#
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	7.975	152		105235	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136		447878	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164		311441	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188		694752	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240		765316	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264		820628	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.387	96		5210	2.210	ng/uL	0.00
4) Pyridine-d5	3.817	84		14188	2.182	ng/ul	0.00
7) Phenol-d5	7.122	99		92044	9.963	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67		64140	11.661	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132		75302	10.805	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113		63719	8.262	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128		40162	12.435	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143		40947	11.564	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165		81669	11.022	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131		80843	7.660	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166		321973	13.161	ng/ul	-0.01
49) Acenaphthylene-d8	14.292	160		336105	11.597	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143		48054	10.831	ng/ul	-0.01
60) Fluorene-d10	15.592	176		279708	13.591	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200		39098	9.500	ng/ul	-0.01
73) Anthracene-d10	17.451	188		451304	13.567	ng/ul	-0.01
81) Pyrene-d10	19.680	212		622040	14.882	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264		617604	14.452	ng/ul	-0.02
Target Compounds							
16) Acetophenone	8.793	105		16065m	1.322	ng/ul	Qvalue

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

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