

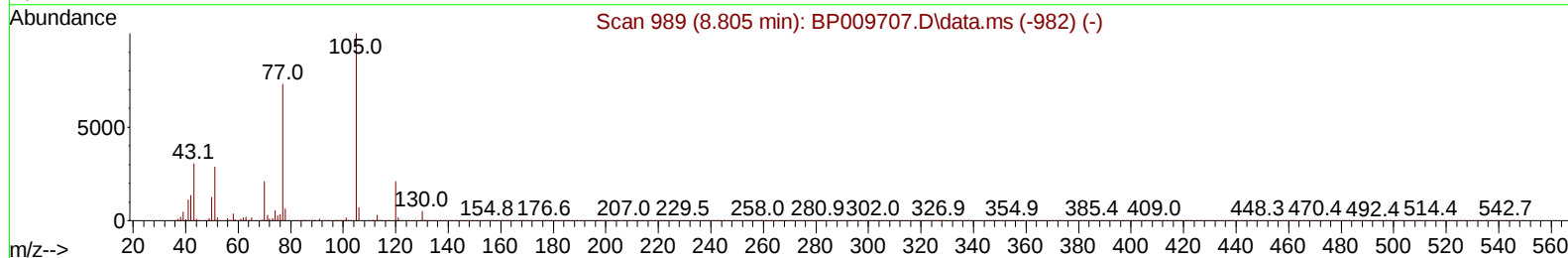
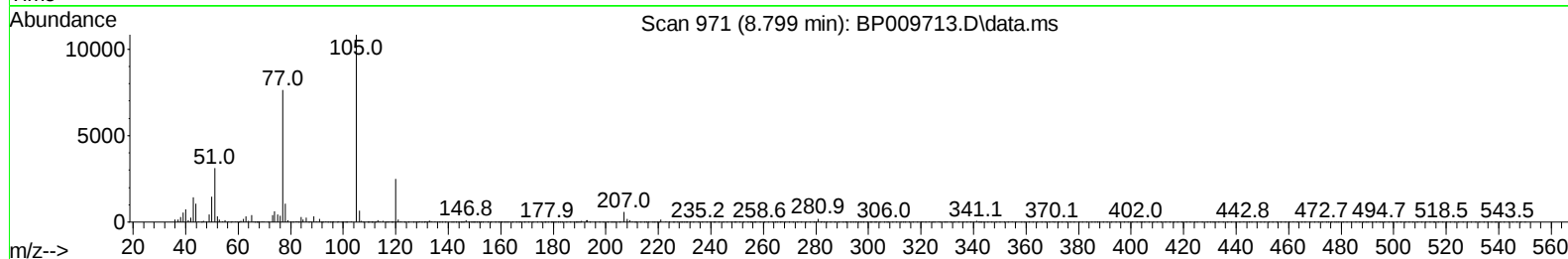
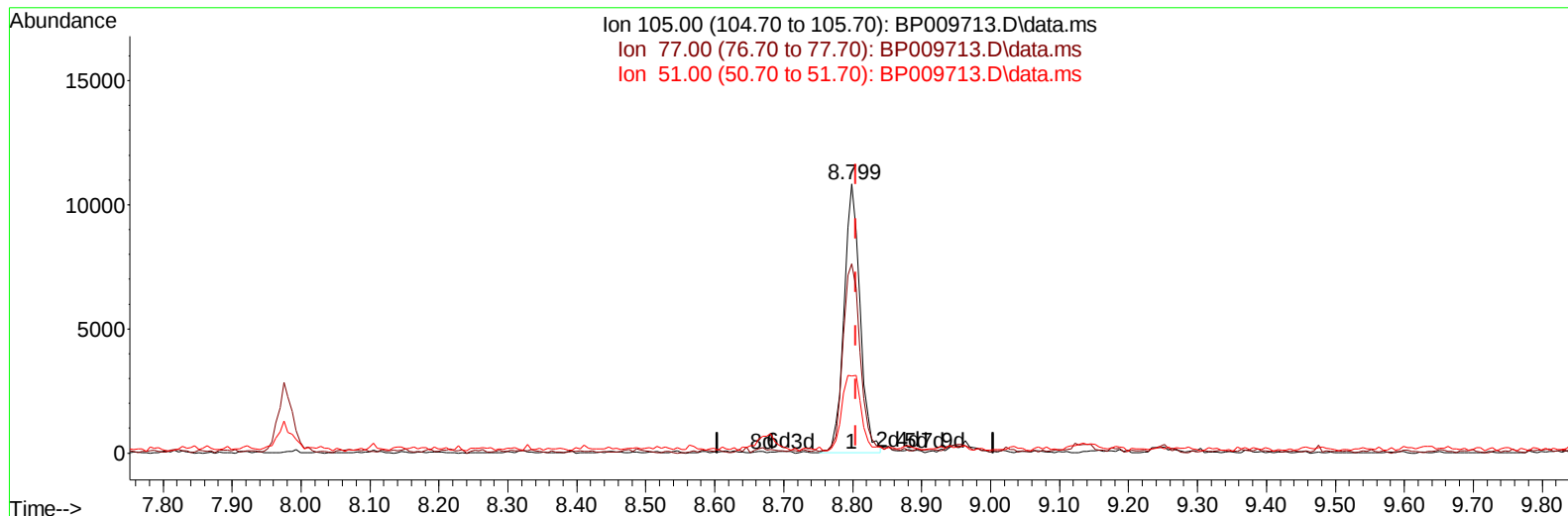
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009713.D
 Acq On : 08 Apr 2022 12:33
 Operator : CG/JU
 Sample : N2183-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R31

Manual IntegrationsAPPROVED

Quant Time: Apr 08 22:26:34 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 04/11/2022
 Supervised By :Yogesh Patel 04/12/2022



TIC: BP009713.D\data.ms

(16) Acetophenone

8.799min (-0.006) 1.45 ng/u1

response 17421

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	70.42
51.00	17.30	28.71#
0.00	0.00	0.00

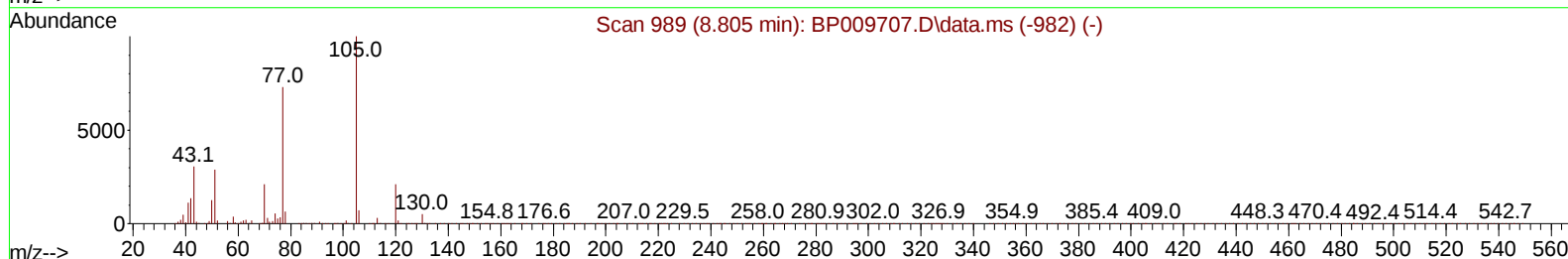
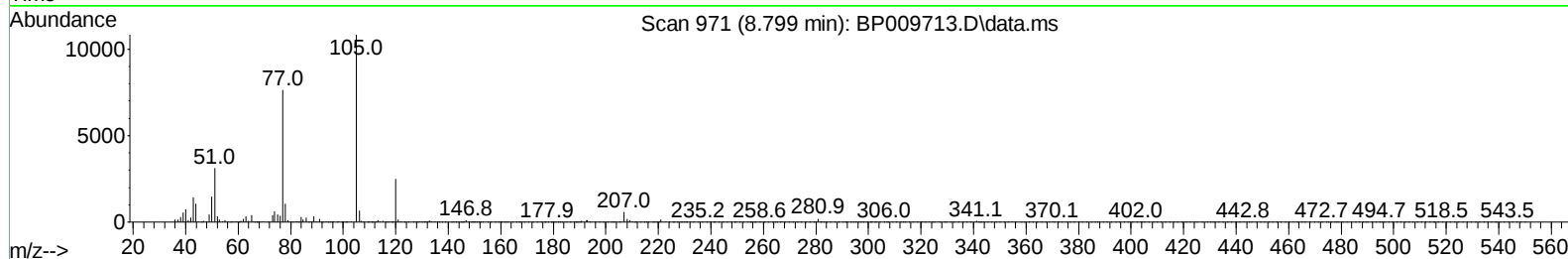
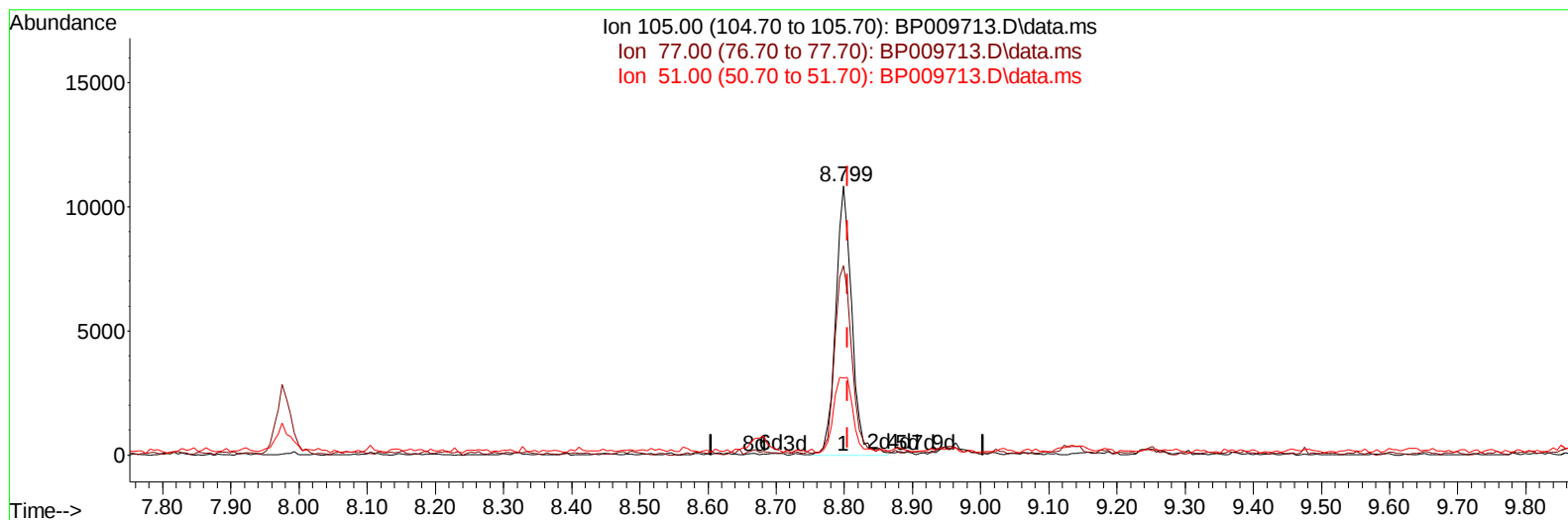
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TIC: BP009713.D\data.ms

(16) Acetophenone

8.799min (-0.006) 1.48 ng/ul m

response 17740

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	70.42
51.00	17.30	28.71#
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	7.975	152	103942	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	465489	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	343962	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	774684	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.439	240	843340	20.000	ng/ul	# 0.00
88) Perylene-d12	23.916	264	837956	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	7246	3.111	ng/uL	0.00
4) Pyridine-d5	3.811	84	57433	8.944	ng/ul	0.00
7) Phenol-d5	7.128	99	123613	13.546	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	81475	14.997	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	100036	14.533	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	107622	14.128	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	50775	15.126	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	52428	14.246	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	104726	13.599	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	142546	12.996	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	430266	15.925	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	458383	14.320	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	70954	14.480	ng/ul	0.00
60) Fluorene-d10	15.598	176	370264	16.290	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	55845	12.169	ng/ul	0.00
73) Anthracene-d10	17.457	188	630816	17.007	ng/ul	0.00
81) Pyrene-d10	19.686	212	940412	20.417	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	903569	20.706	ng/ul	0.00
Target Compounds						
16) Acetophenone	8.799	105	17740m	1.478	ng/ul	Qvalue

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

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