

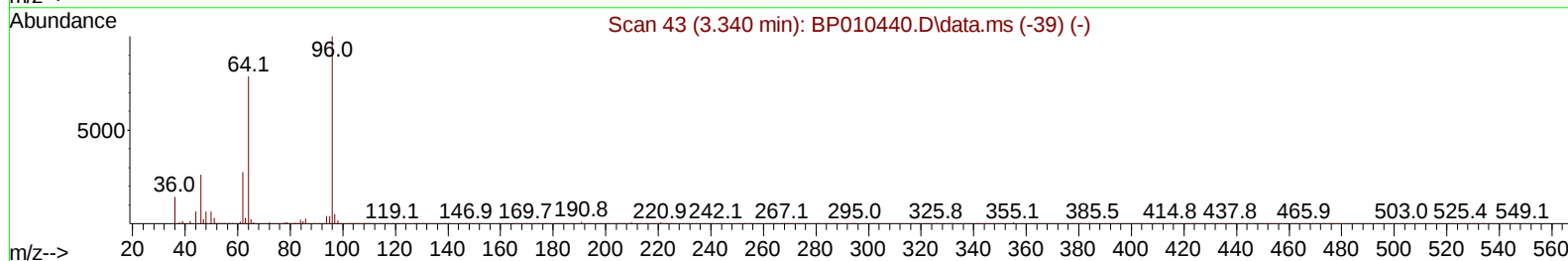
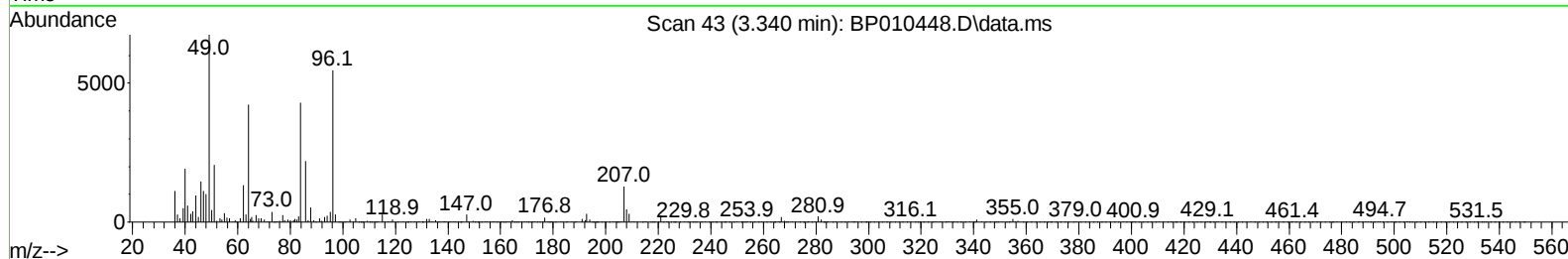
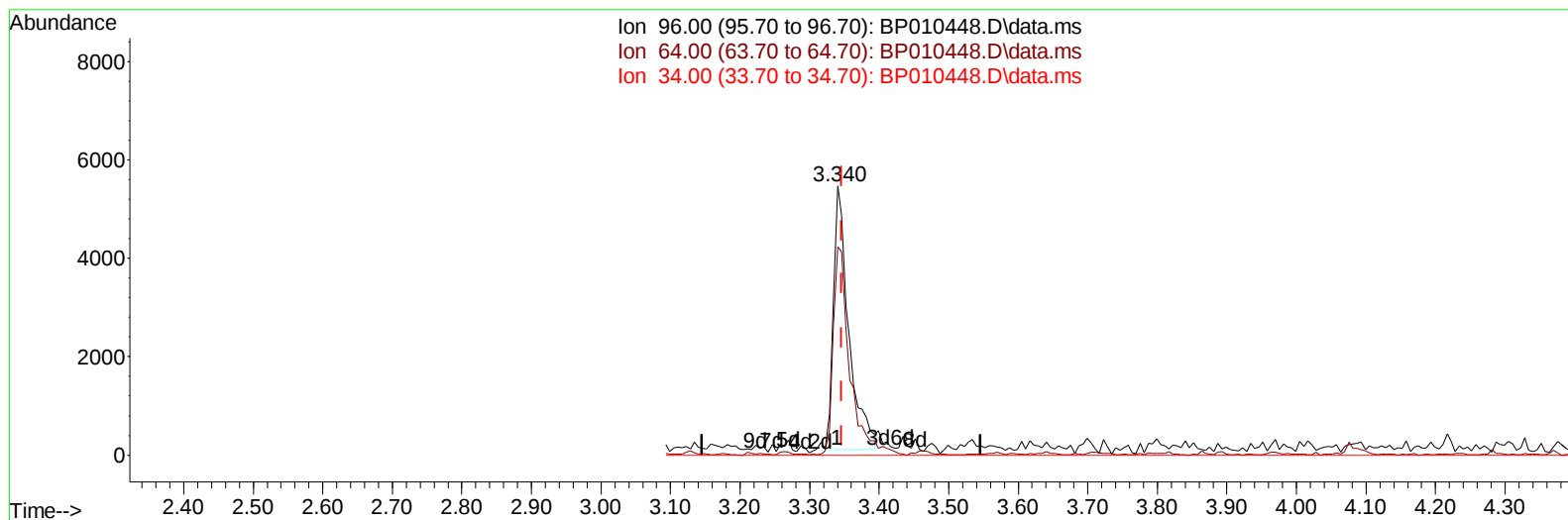
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
Data File : BP010448.D
Acq On : 20 May 2022 22:33
Operator : CG/JU
Sample : N2843-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4J3

Manual IntegrationsAPPROVED

Quant Time: May 21 01:20:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
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Reviewed By :Jagrut Upadhyay 05/23/2022
Supervised By :mohammad ahmed 05/25/2022



TIC: BP010448.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.340min (-0.006) 2.75 ng/uL

response 8219

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	77.52#
34.00	0.00	0.00
0.00	0.00	0.00

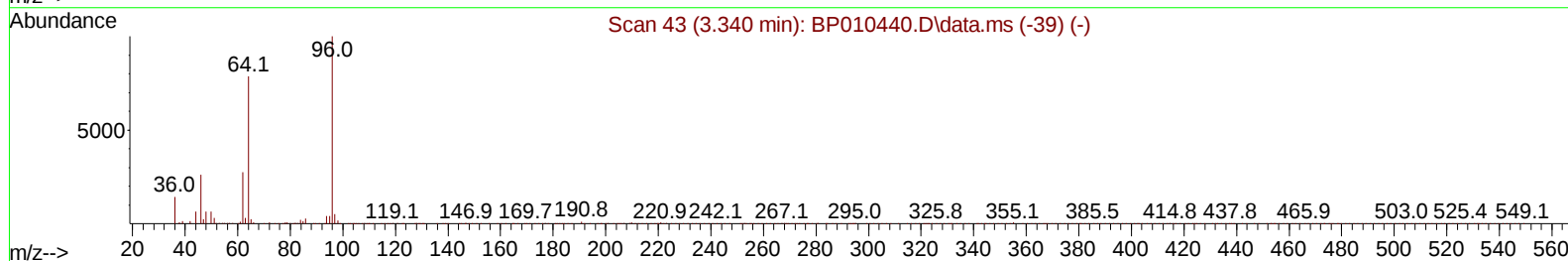
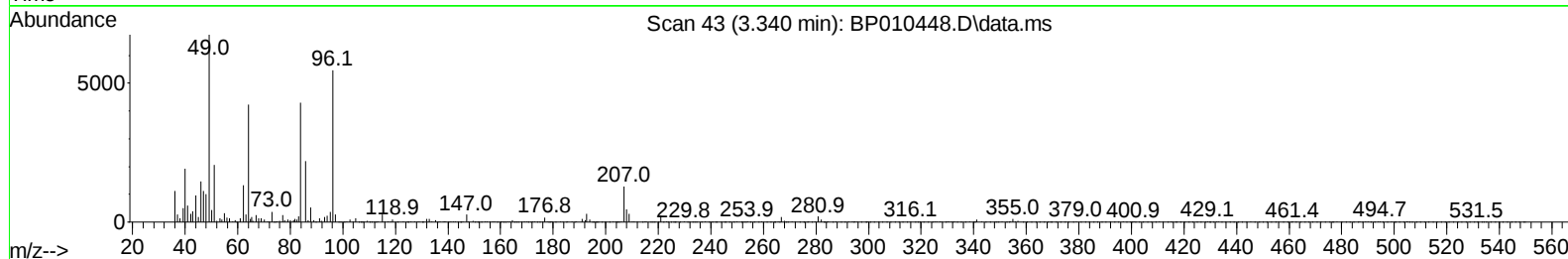
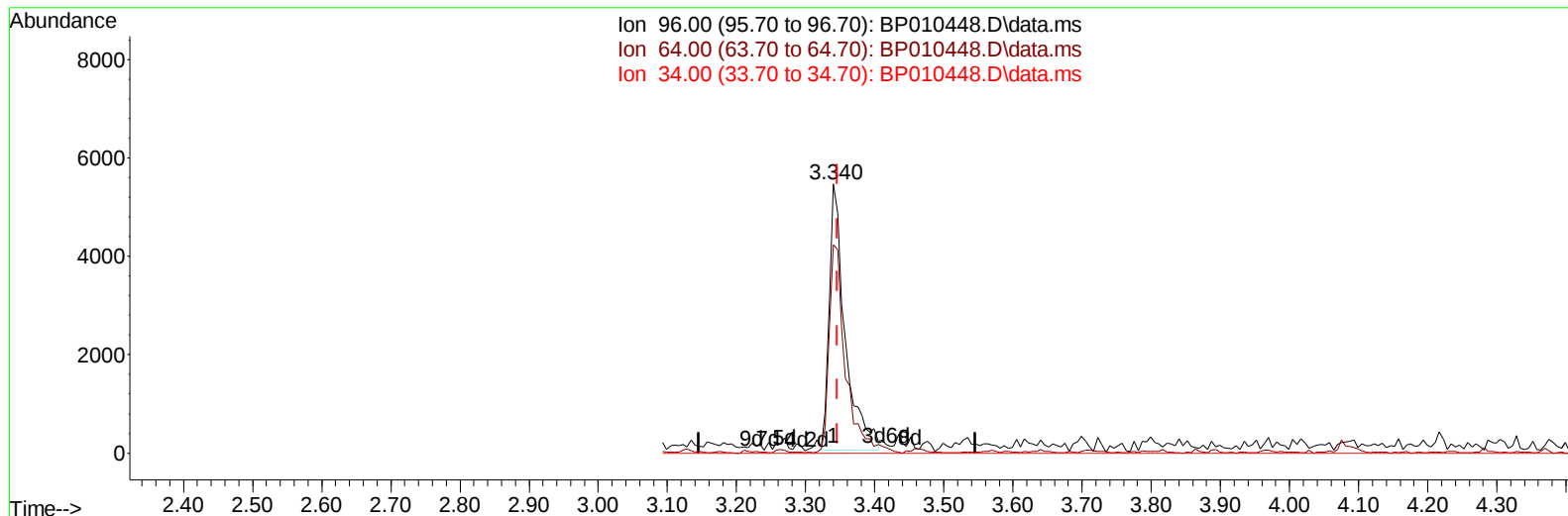
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TIC: BP010448.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.340min (-0.006) 2.88 ng/uL m

response 8629

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	77.52#
34.00	0.00	0.00
0.00	0.00	0.00

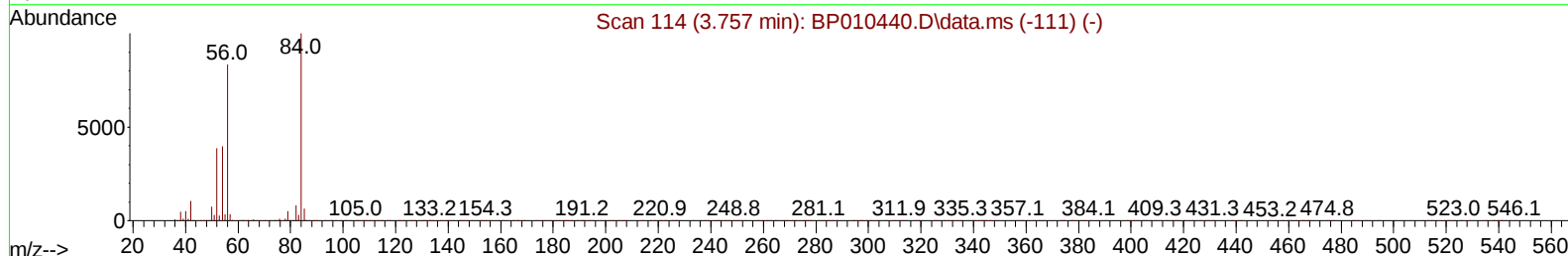
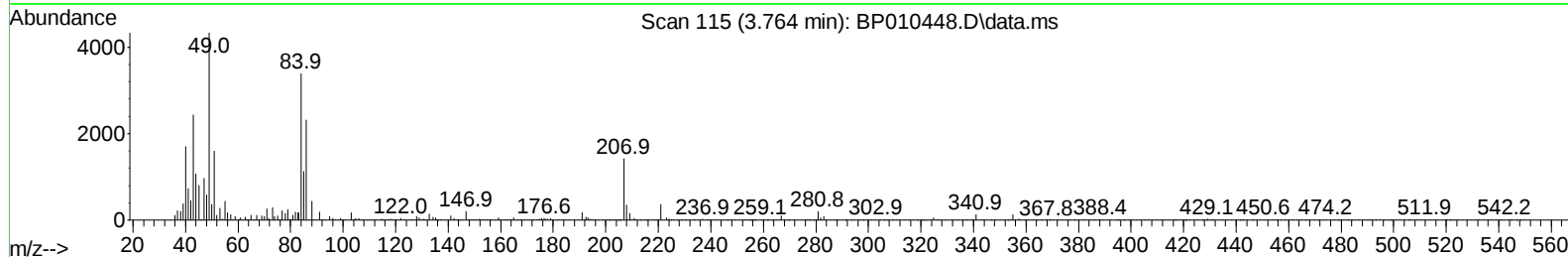
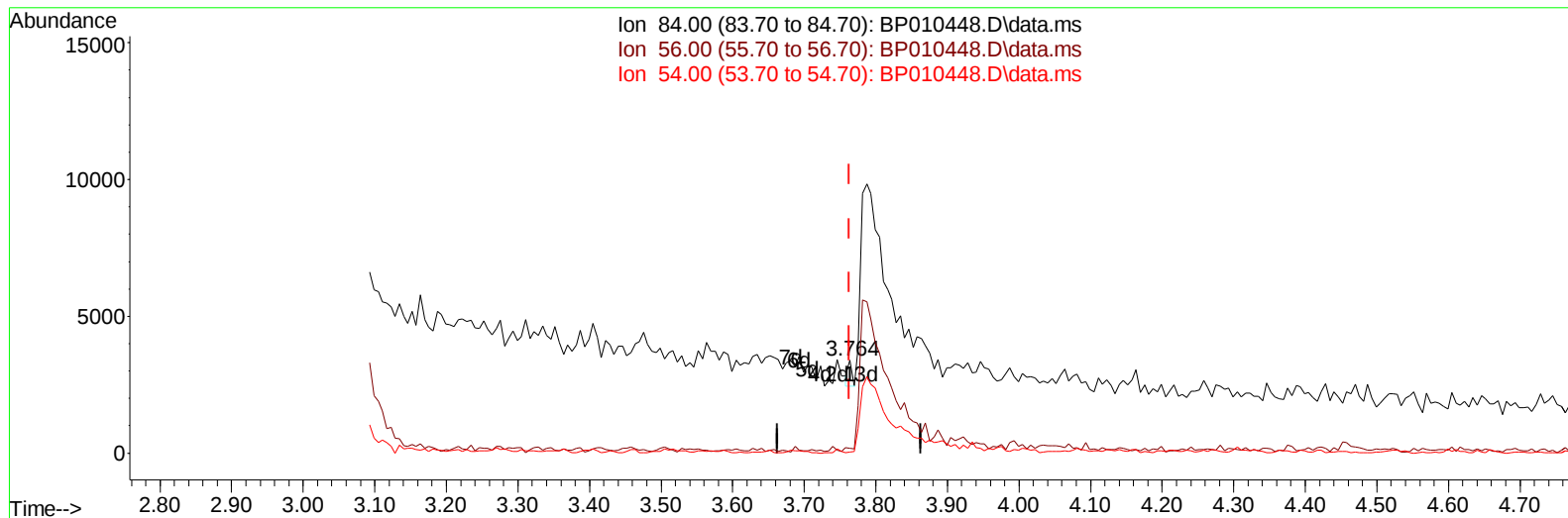
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TIC: BP010448.D\data.ms

(4) Pyridine-d5 (S)

3.764min (+ 0.000) 0.04 ng/u1

response 332

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	5.04#
54.00	24.80	1.06#
0.00	0.00	0.00

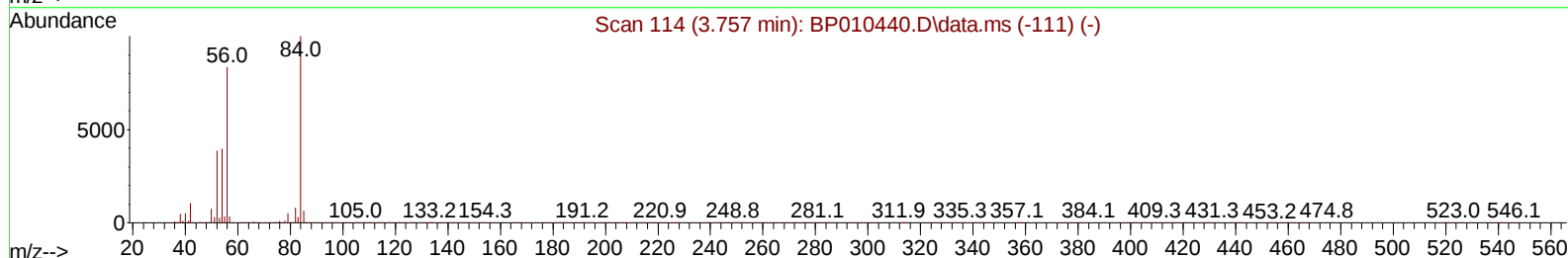
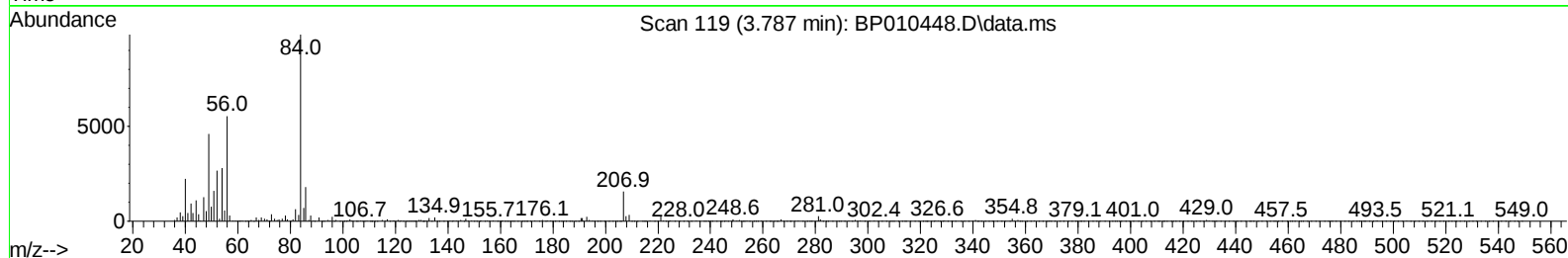
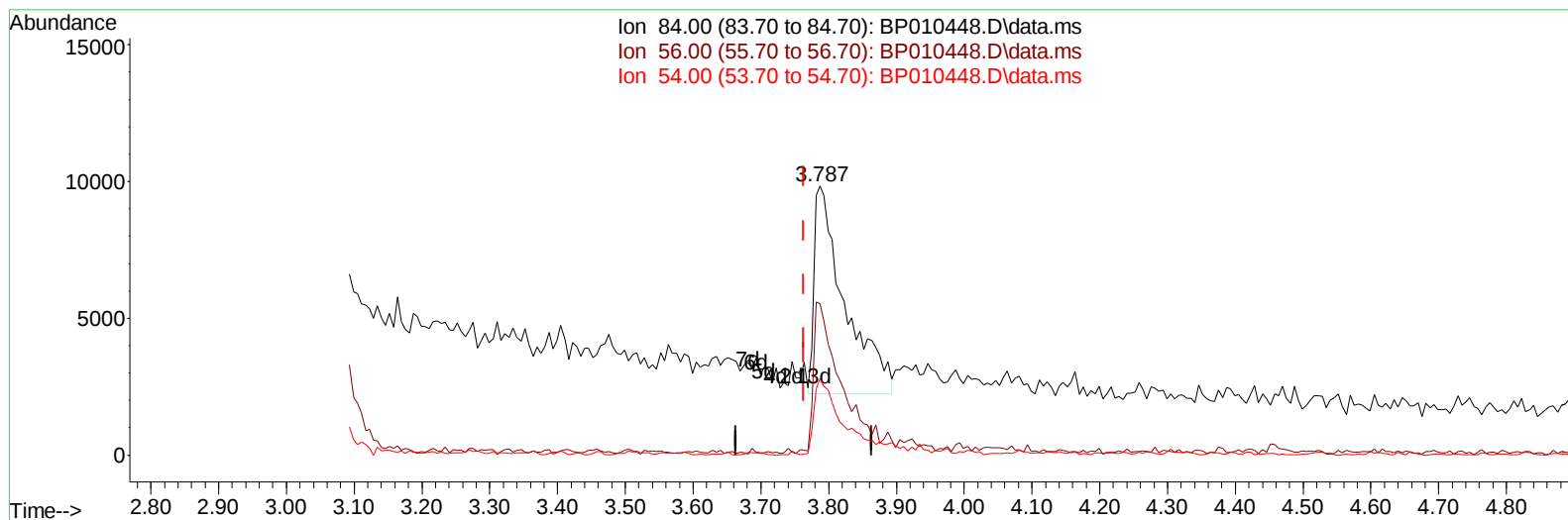
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Instrument :
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TIC: BP010448.D\data.ms

(4) Pyridine-d5 (S)

3.787min (+ 0.024) 2.79 ng/u1 m

response 23699

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	56.28#
54.00	24.80	28.33
0.00	0.00	0.00

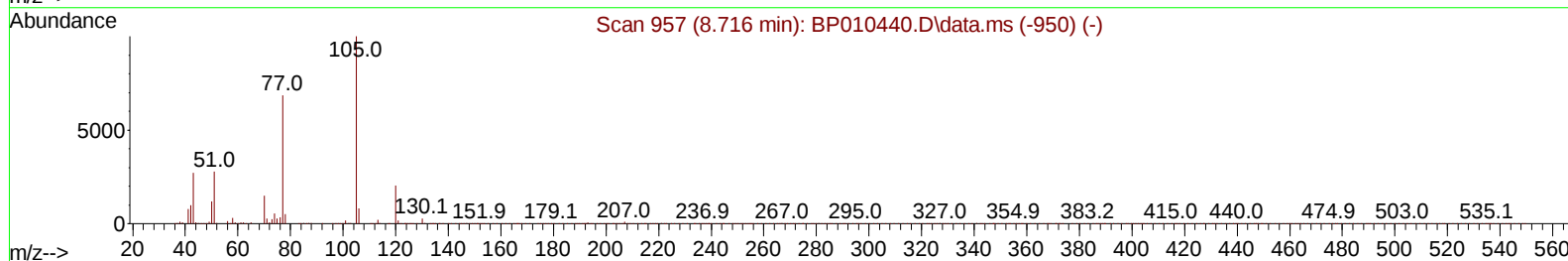
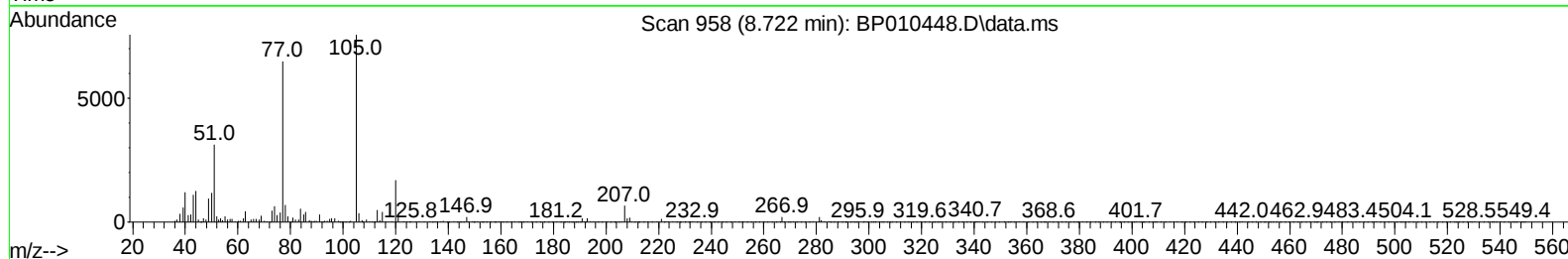
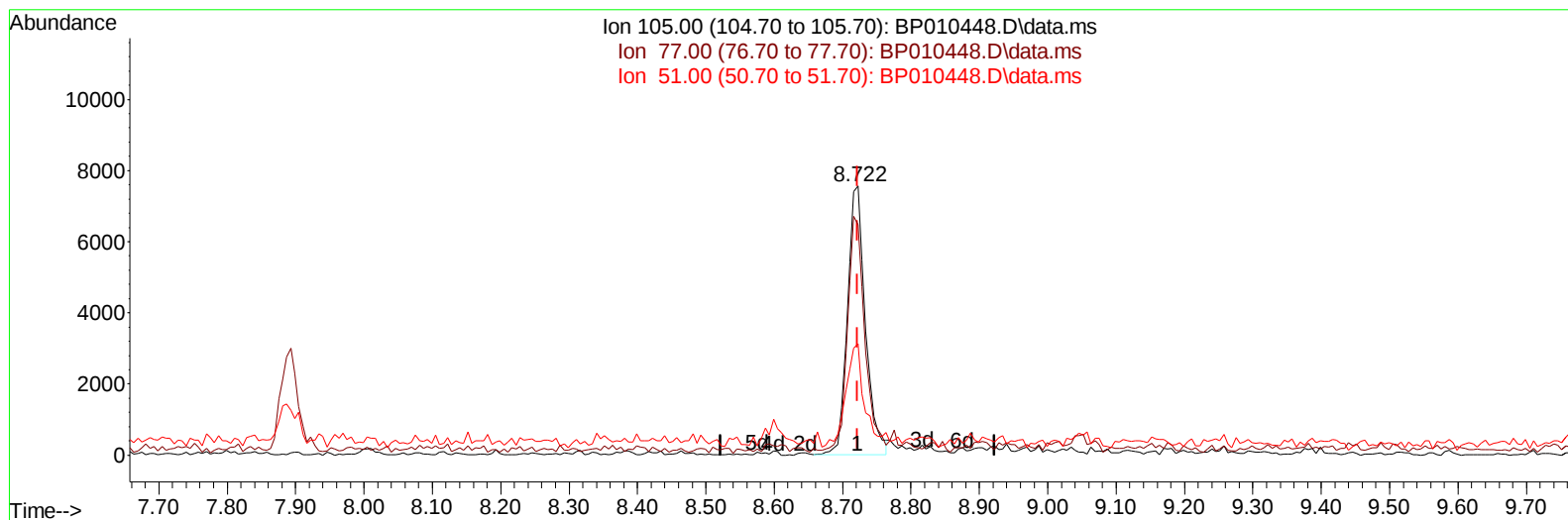
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TIC: BP010448.D\data.ms

(16) Acetophenone

8.722min (+ 0.000) 0.97 ng/u1

response 13799

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	85.65#
51.00	17.30	41.40#
0.00	0.00	0.00

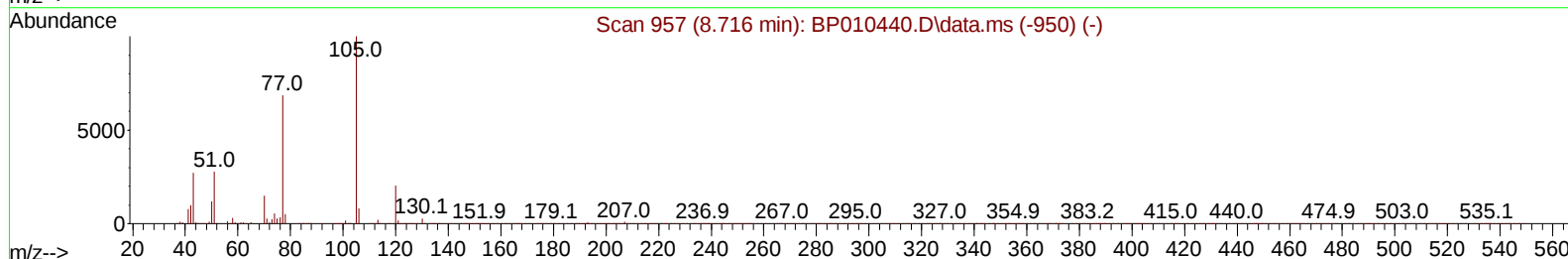
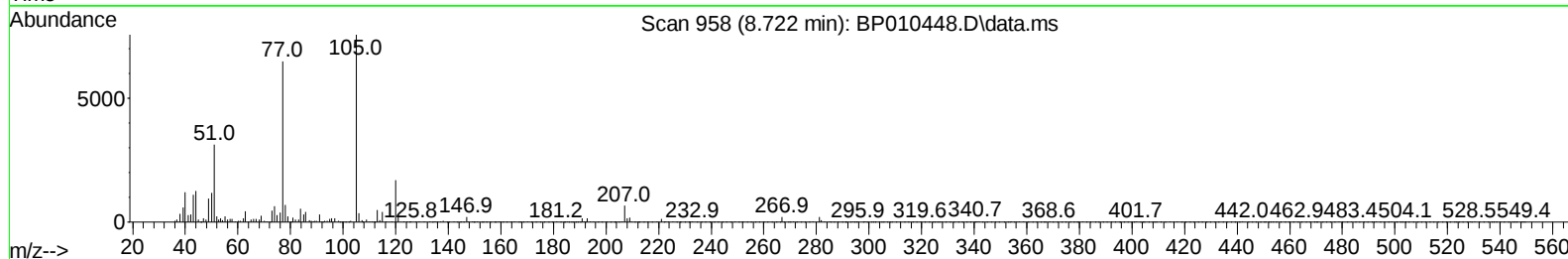
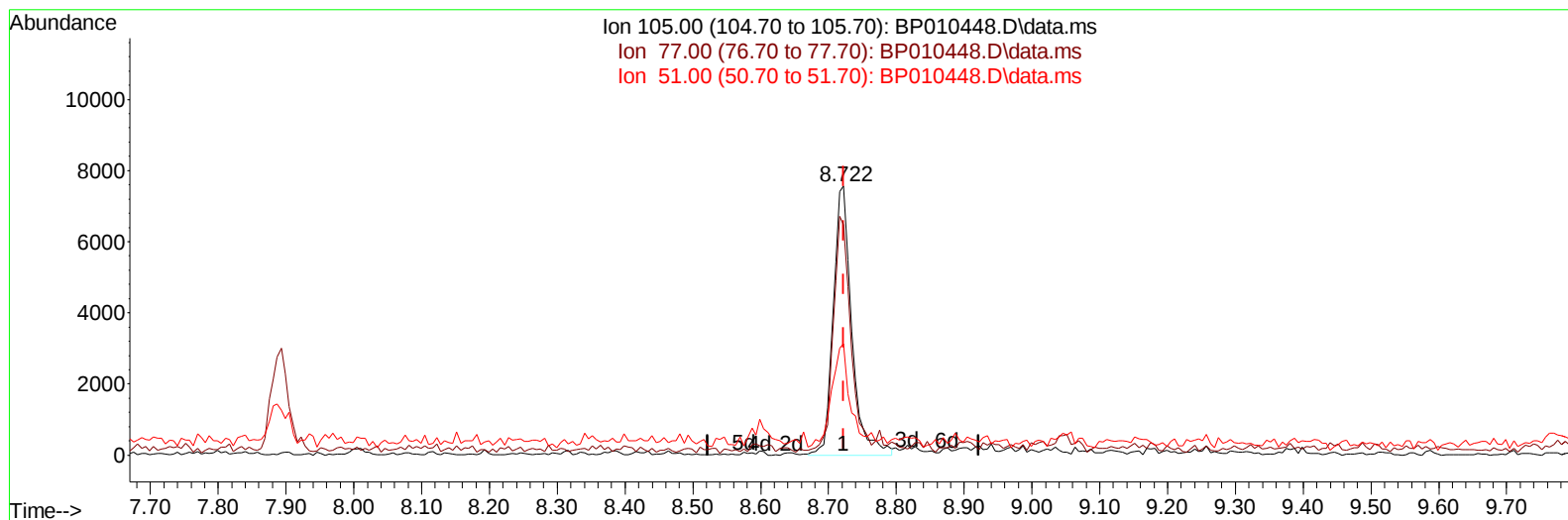
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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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TIC: BP010448.D\data.ms

(16) Acetophenone

8.722min (+ 0.000) 1.00 ng/ul m

response 14311

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	85.65#
51.00	17.30	41.40#
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	7.893	152		127553	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136		565881	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164		389517	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188		835383	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240		840090	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264		792982	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.340	96		8629m	2.884	ng/uL	0.00
4) Pyridine-d5	3.787	84		23699m	2.789	ng/ul	0.02
7) Phenol-d5	7.058	99		173259	16.184	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67		124227	17.271	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132		141376	17.258	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113		121726	13.379	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128		77295	17.649	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143		74533	16.059	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165		136994	15.610	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131		168060	12.879	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166		498486	17.370	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160		586853	17.728	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143		51298	10.043	ng/ul	0.00
60) Fluorene-d10	15.516	176		458613	18.369	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200		42908	8.782	ng/ul	0.00
73) Anthracene-d10	17.375	188		738216	19.515	ng/ul	0.00
81) Pyrene-d10	19.604	212		987229	22.128	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.621	264		840908	21.118	ng/ul	-0.01
Target Compounds							
16) Acetophenone	8.722	105		14311m	1.001	ng/ul	Qvalue

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

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