

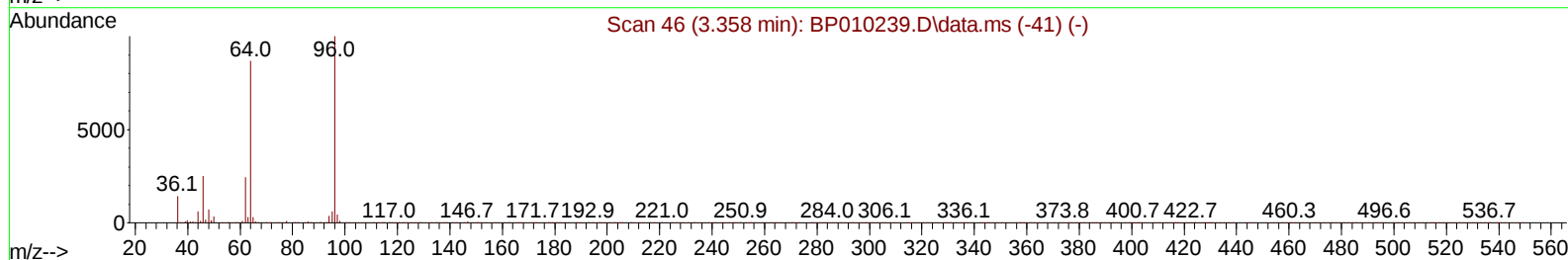
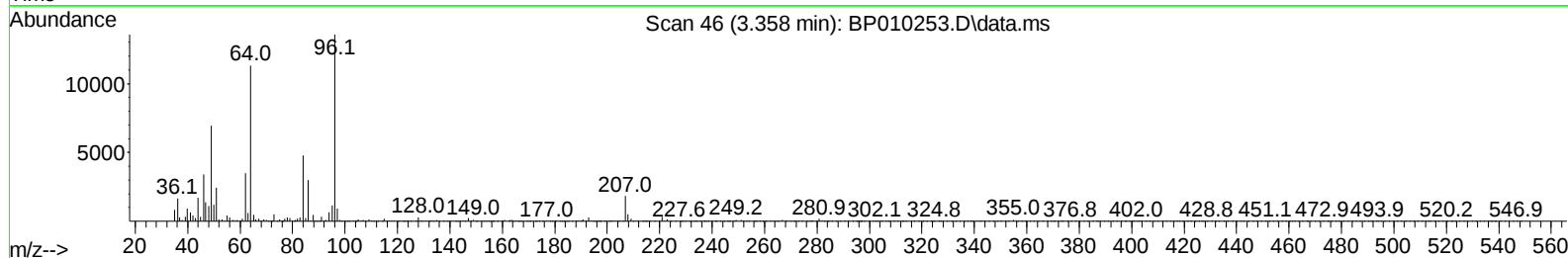
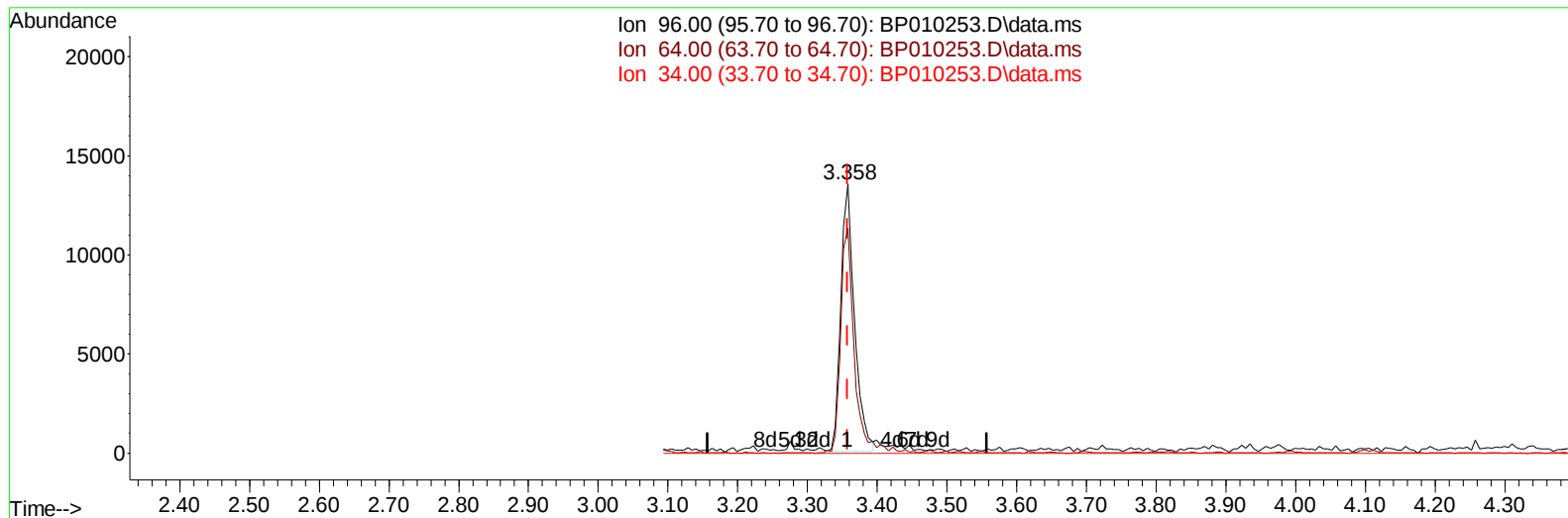
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010253.D
 Acq On : 05 May 2022 18:33
 Operator : CG/JU
 Sample : N2678-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0B7

Manual IntegrationsAPPROVED

Quant Time: May 06 02:07:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 06 01:05:07 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/10/2022
 Supervised By :mohammad ahmed 05/10/2022



TIC: BP010253.D\data.ms

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

3.358min (+ 0.000) 4.37 ng/uL

response 18337

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.50#
34.00	0.00	0.00
0.00	0.00	0.00

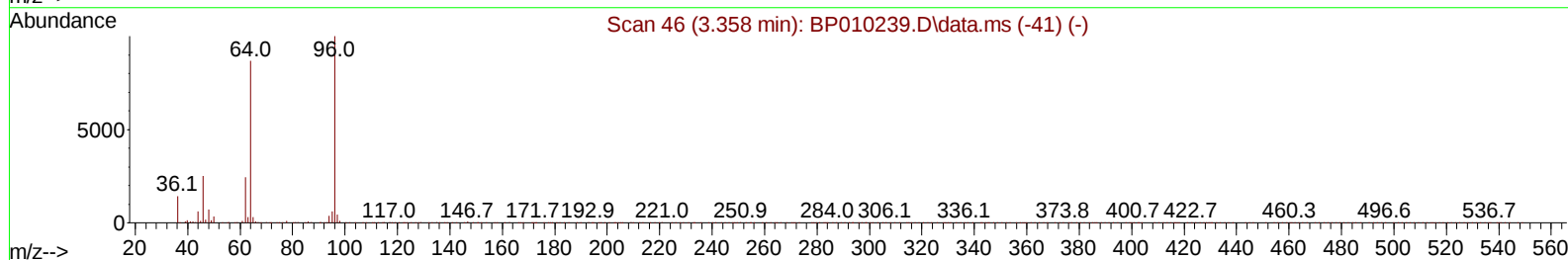
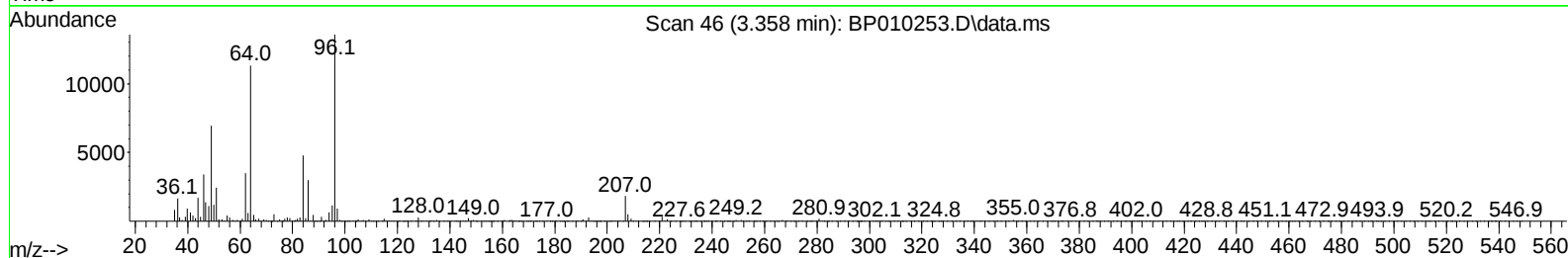
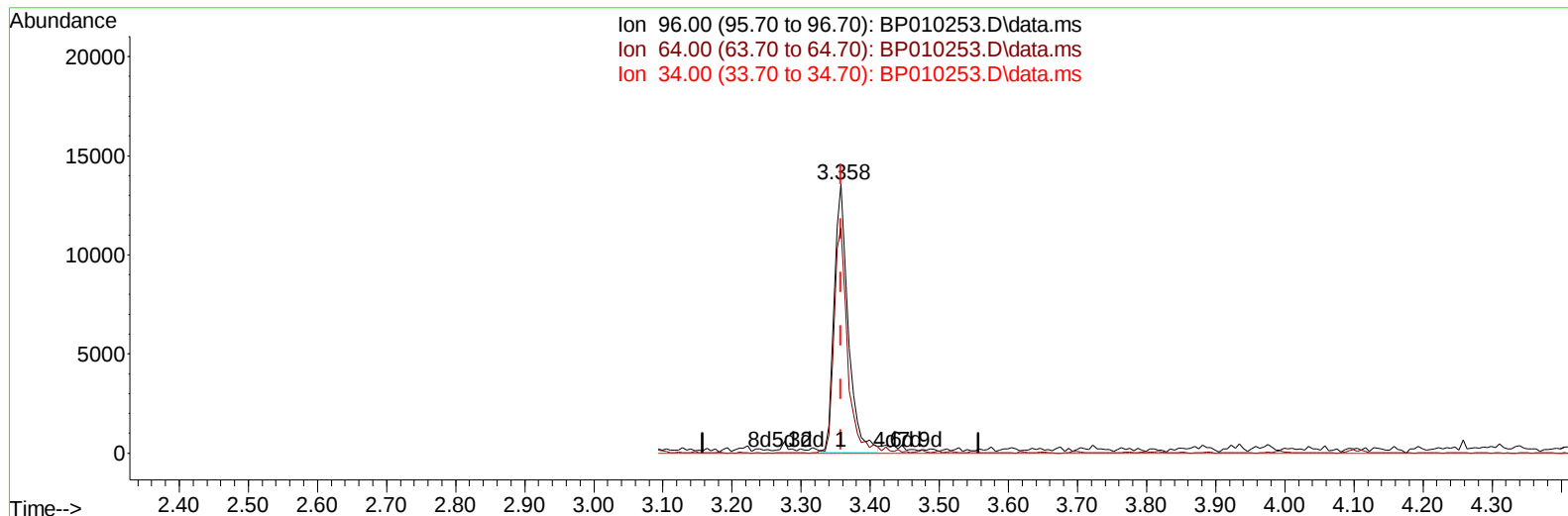
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010253.D
 Acq On : 05 May 2022 18:33
 Operator : CG/JU
 Sample : N2678-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 BH0B7

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 Supervised By :mohammad ahmed 05/10/2022



TIC: BP010253.D\data.ms

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

3.358min (+ 0.000) 4.50 ng/uL m

response 18885

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.50#
34.00	0.00	0.00
0.00	0.00	0.00

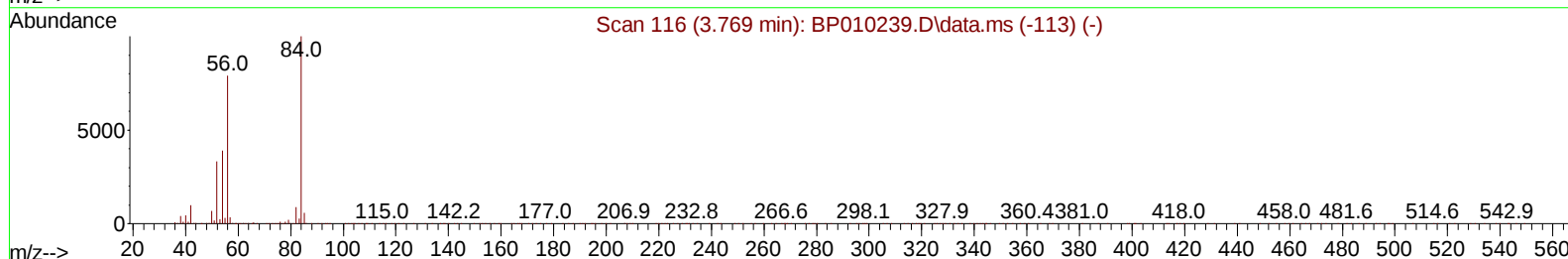
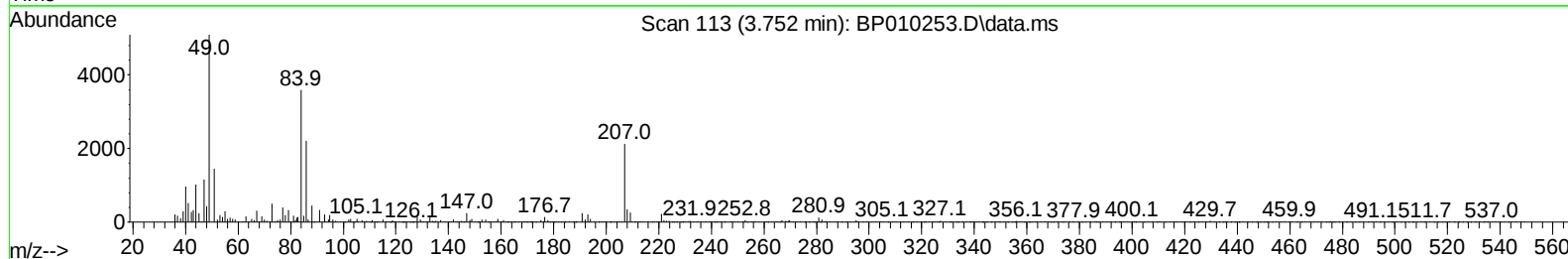
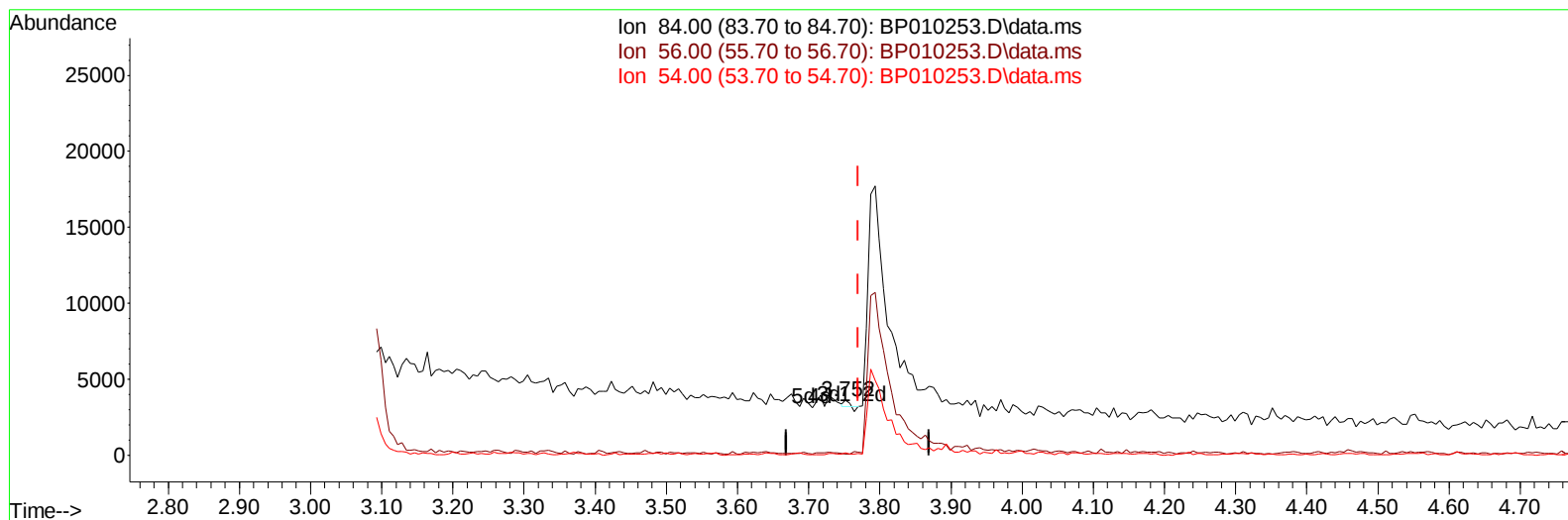
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010253.D
 Acq On : 05 May 2022 18:33
 Operator : CG/JU
 Sample : N2678-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP010253.D\data.ms

(4) Pyridine-d5 (S)

3.752min (-0.018) 0.01 ng/u1

response 70

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	2.20#
54.00	24.80	3.84#
0.00	0.00	0.00

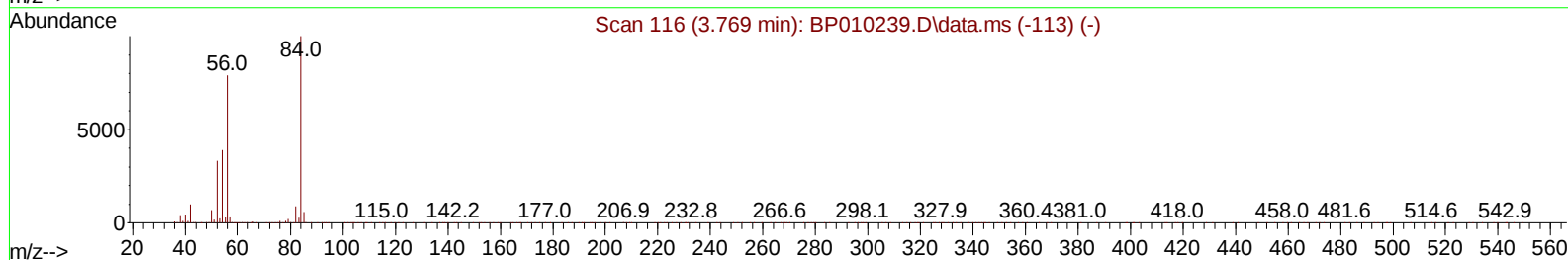
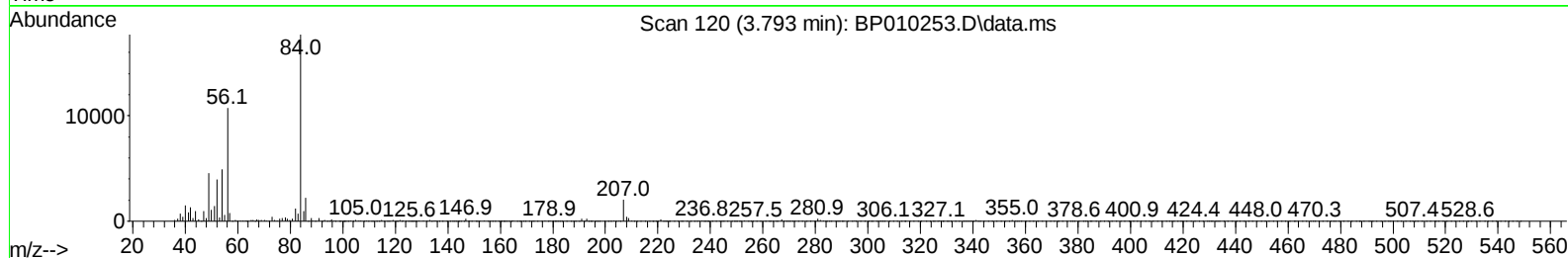
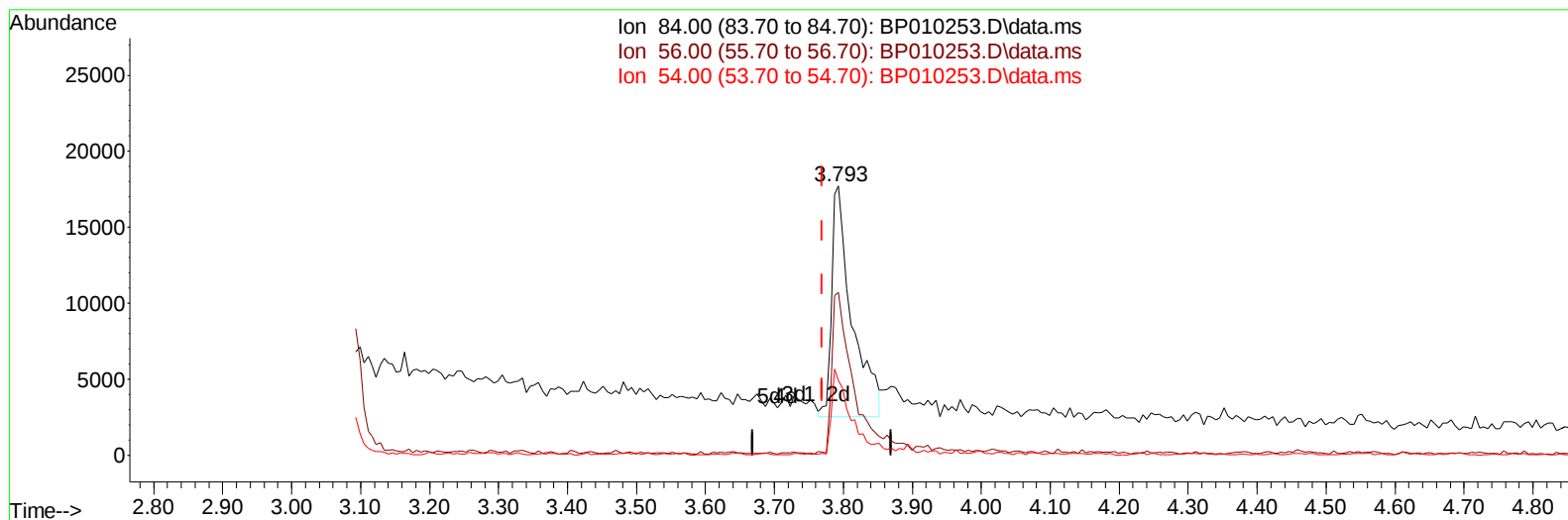
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010253.D
 Acq On : 05 May 2022 18:33
 Operator : CG/JU
 Sample : N2678-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP010253.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.024) 2.62 ng/ul m

response 30903

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	60.54#
54.00	24.80	27.78
0.00	0.00	0.00

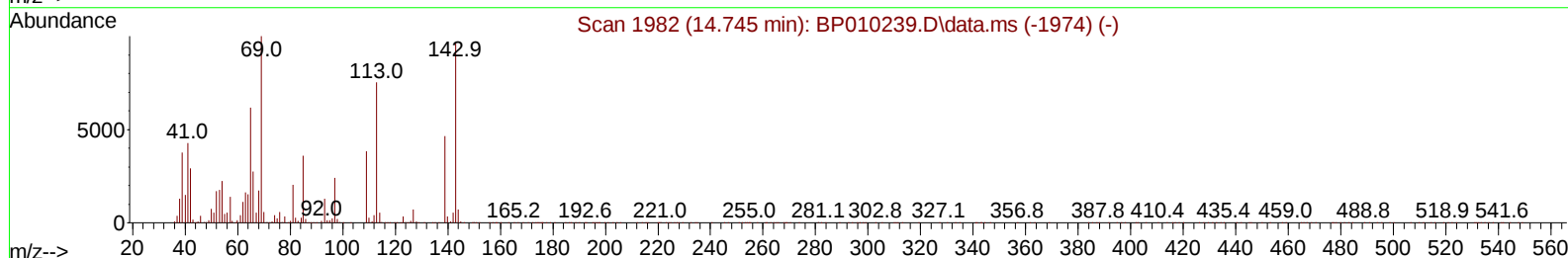
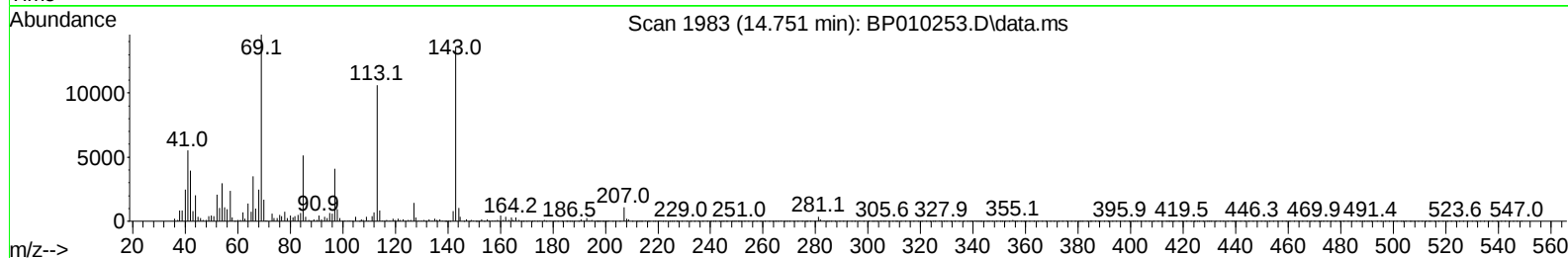
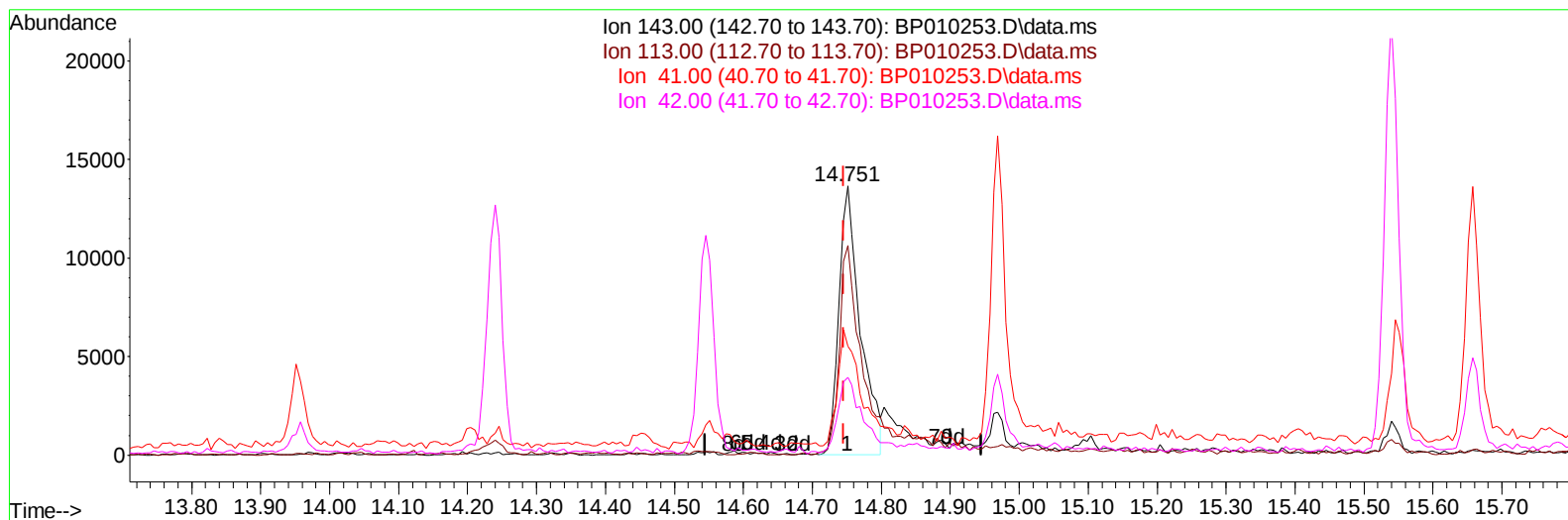
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
Data File : BP010253.D
Acq On : 05 May 2022 18:33
Operator : CG/JU
Sample : N2678-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0B7

Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 05/10/2022

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TIC: BP010253.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 3.98 ng/u1

response 29356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.92#
41.00	23.20	40.50#
42.00	18.00	28.92#

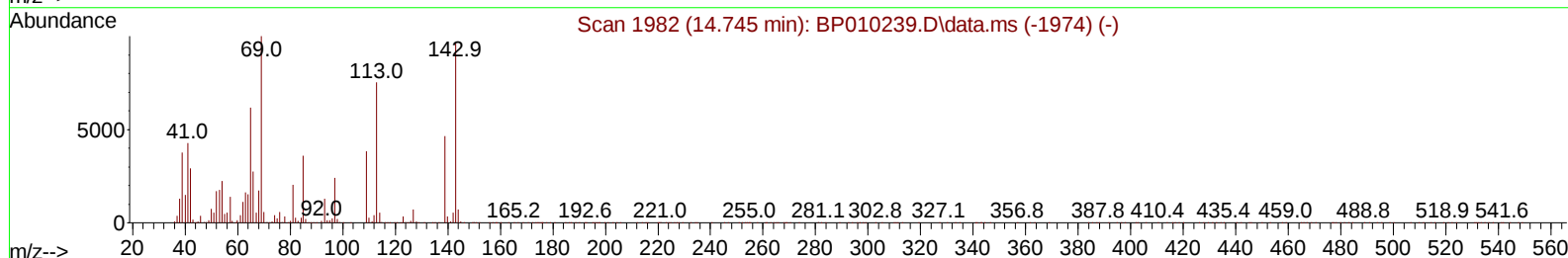
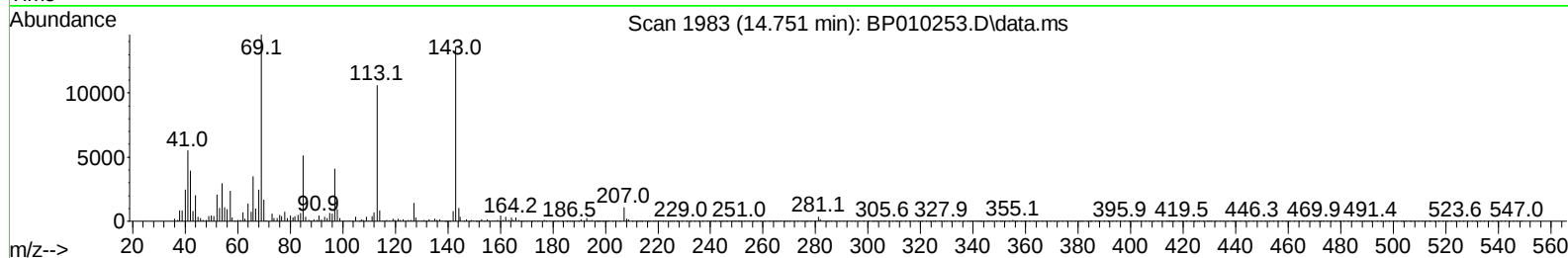
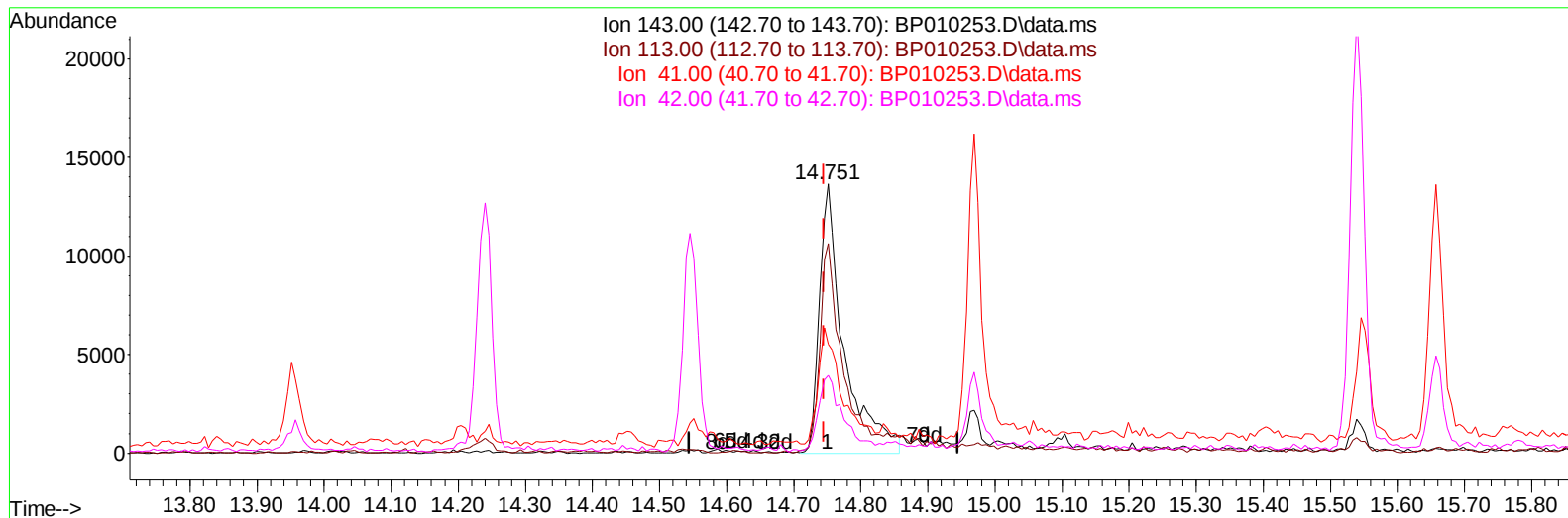
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
Data File : BP010253.D
Acq On : 05 May 2022 18:33
Operator : CG/JU
Sample : N2678-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0B7

Manual IntegrationsAPPROVED

Quant Time: May 06 02:07:35 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
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Reviewed By :Jagrut Upadhyay 05/10/2022
Supervised By :mohammad ahmed 05/10/2022



TIC: BP010253.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.67 ng/ul m

response 34479

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.92#
41.00	23.20	40.50#
42.00	18.00	28.92#

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 Misc :
 ALS Vial : 16 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	183012	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	796498	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.545	164	515197	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	996948	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	646849	20.000	ng/ul	# 0.00
88) Perylene-d12	23.821	264	592135	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	18885m	4.500	ng/uL	0.00
4) Pyridine-d5	3.793	84	30903m	2.617	ng/ul	0.02
7) Phenol-d5	7.081	99	105262	6.323	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	319541	31.990	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	306237	24.840	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	217227	15.750	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	204516	32.284	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	216632	31.312	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	373422	27.725	ng/ul	0.00
31) 4-Chloroaniline-d4	10.852	131	225139	11.658	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1487558	37.118	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1613237	33.260	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	34479m	4.673	ng/ul	0.00
60) Fluorene-d10	15.540	176	1219435	35.373	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	198217	31.444	ng/ul	0.00
73) Anthracene-d10	17.398	188	1822012	37.776	ng/ul	0.00
81) Pyrene-d10	19.628	212	1809817	49.198	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.663	264	1282234	41.305	ng/ul	0.00

Target Compounds Qvalue

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

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