

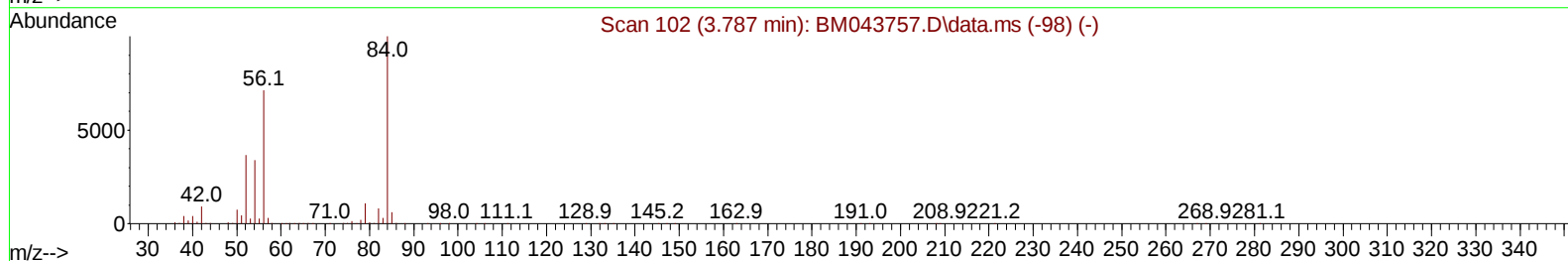
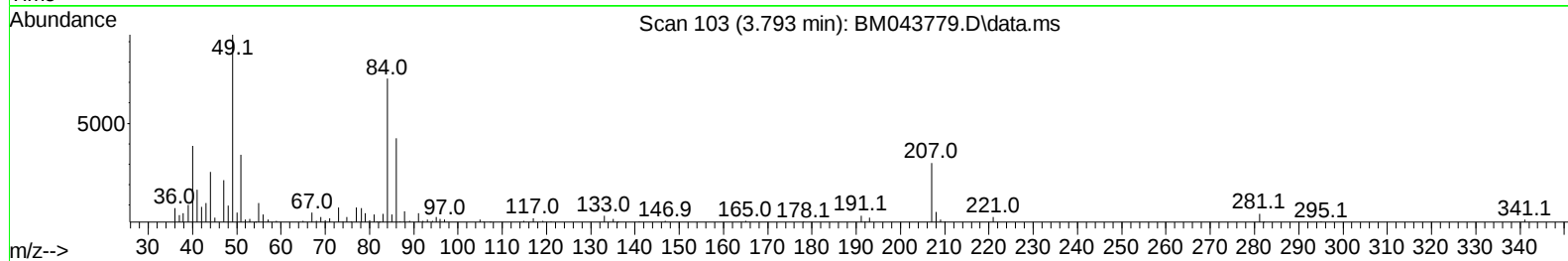
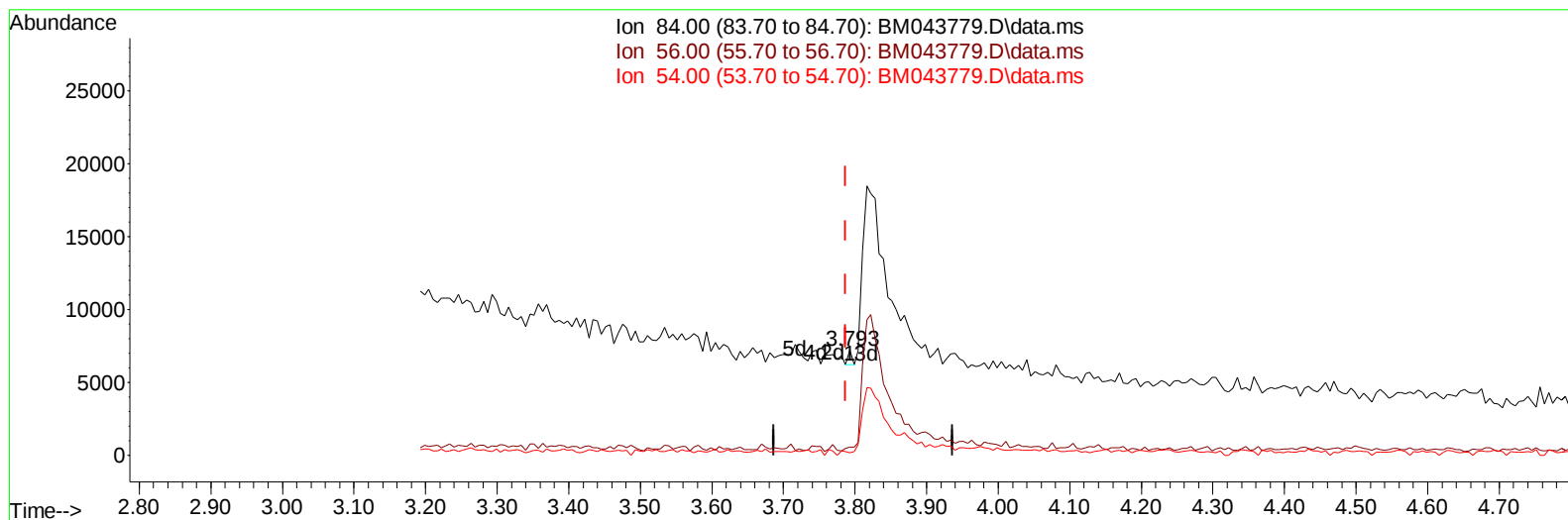
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043779.D
 Acq On : 06 Jan 2024 03:07
 Operator : MA/JU
 Sample : 05952-03MEDL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF88MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 06 03:44:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/08/2024
 Supervised By :mohammad ahmed 01/08/2024



TIC: BM043779.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.006) 0.01 ng/u1

response 354

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.70	7.38#
54.00	35.00	2.08#
0.00	0.00	0.00

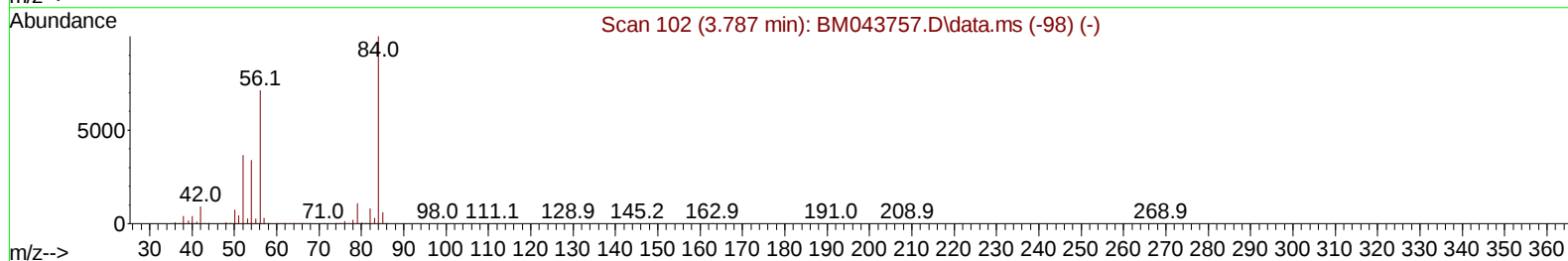
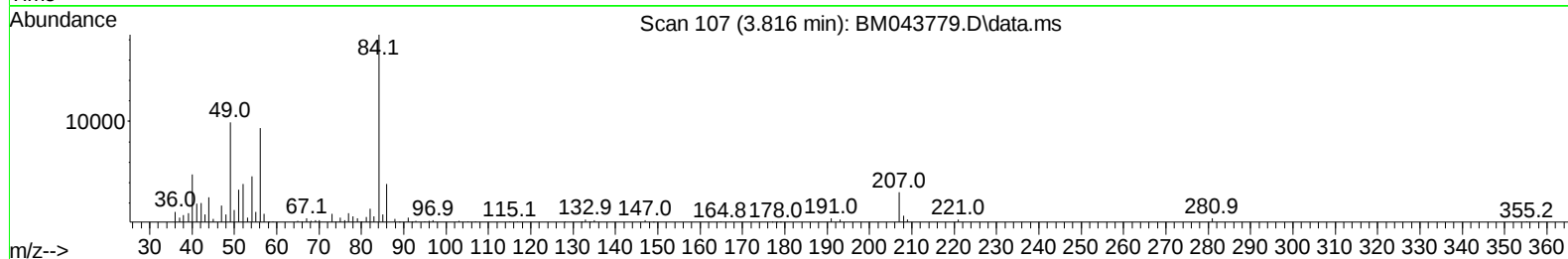
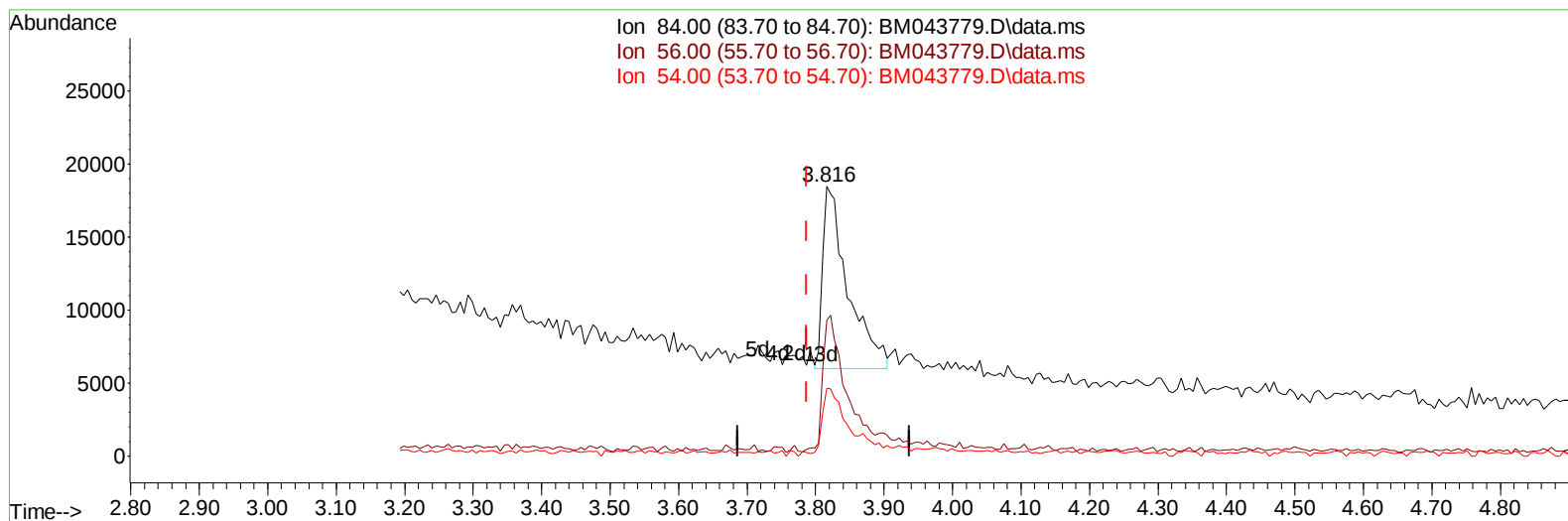
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TIC: BM043779.D\data.ms

(4) Pyridine-d5 (S)

3.816min (+ 0.029) 1.09 ng/u1 m

response 31990

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.70	50.47#
54.00	35.00	25.14#
0.00	0.00	0.00

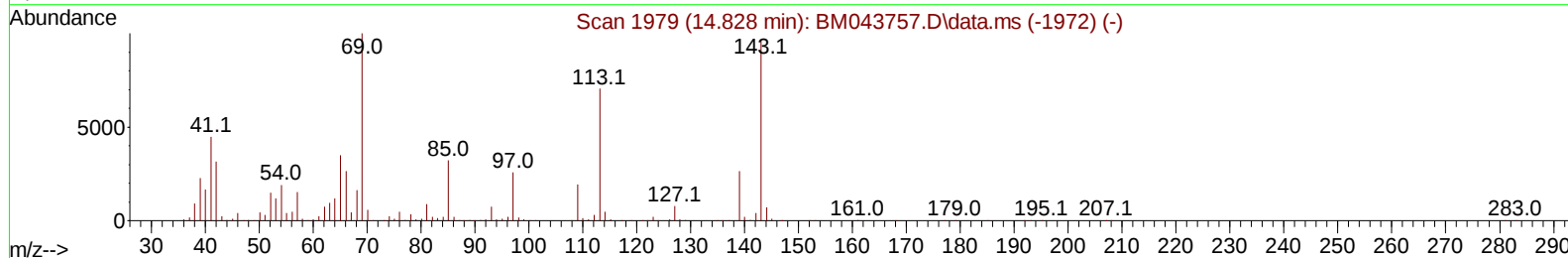
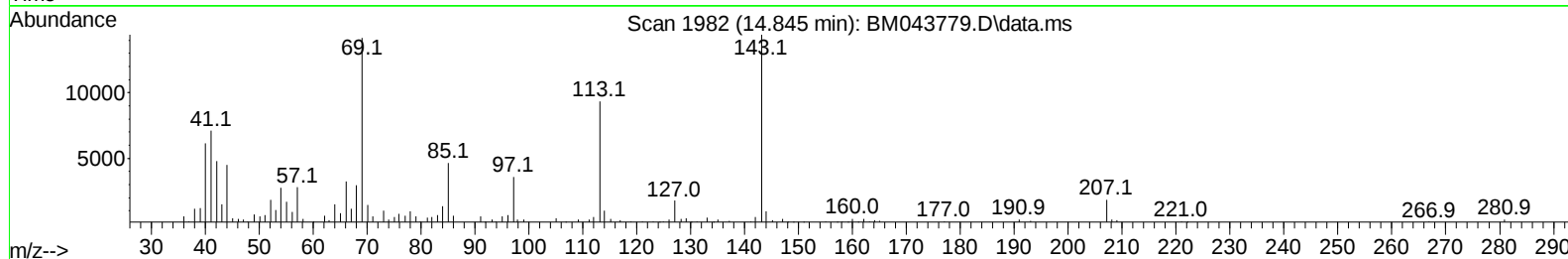
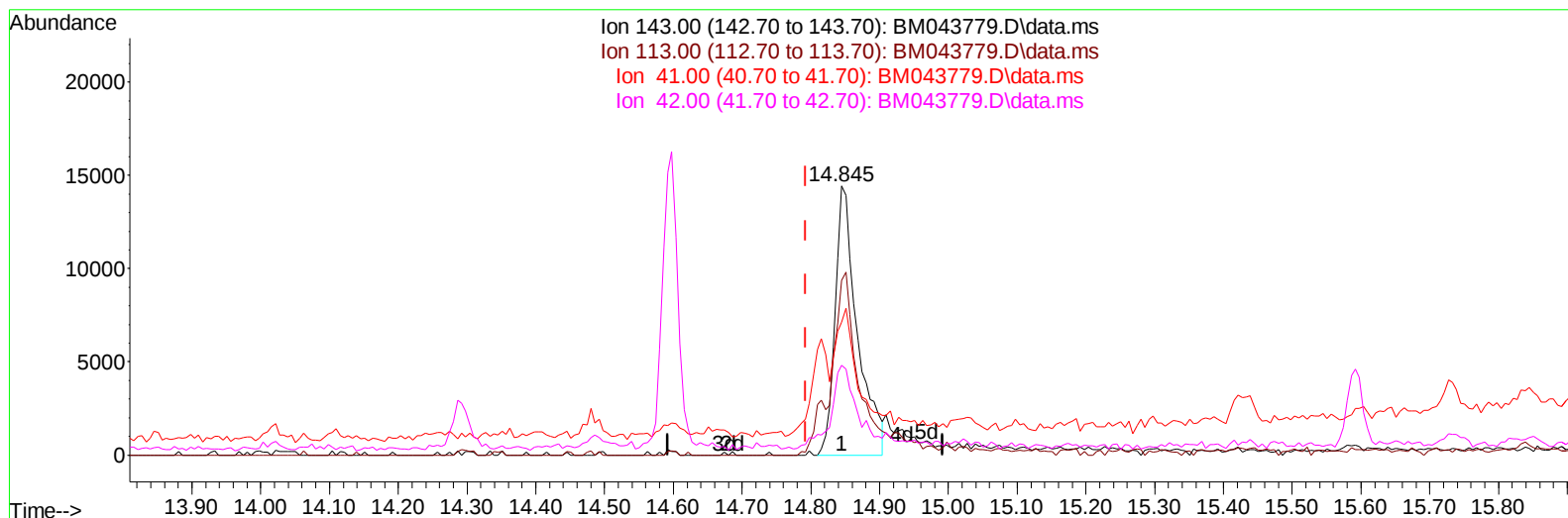
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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BM043779.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (+ 0.053) 2.58 ng/u1

response 32298

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	64.88
41.00	49.10	49.43
42.00	33.20	33.37

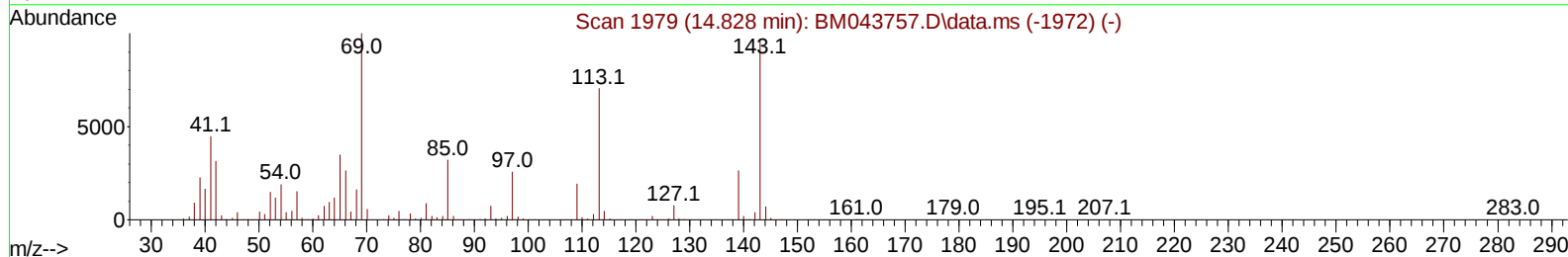
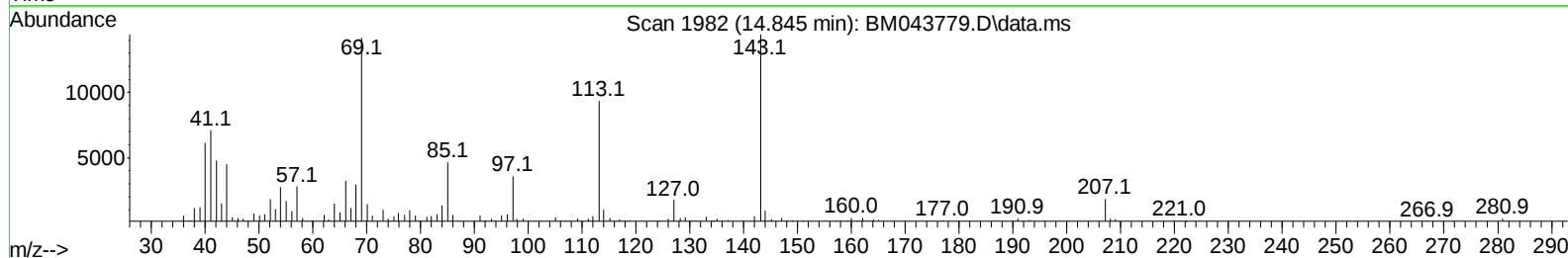
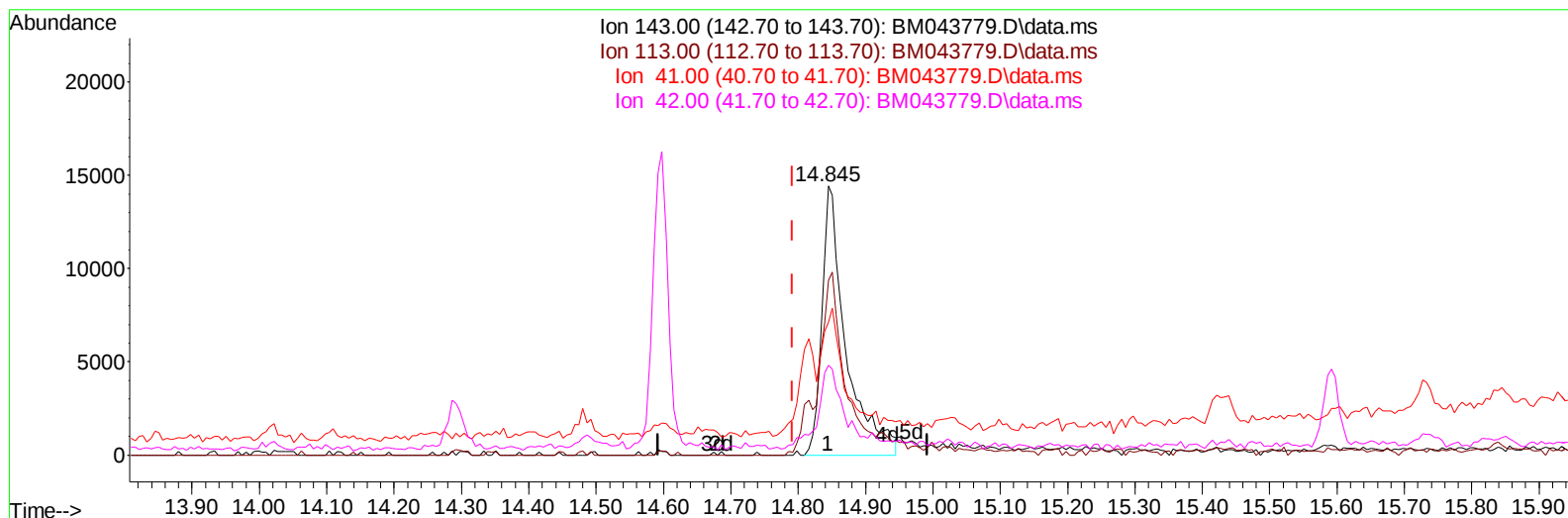
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Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM043779.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (+ 0.053) 2.84 ng/ul m

response 35595

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	64.88
41.00	49.10	49.43
42.00	33.20	33.37

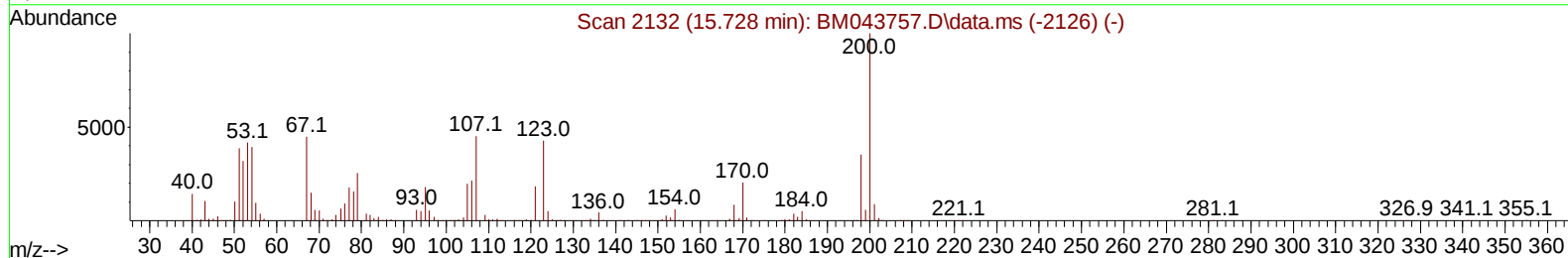
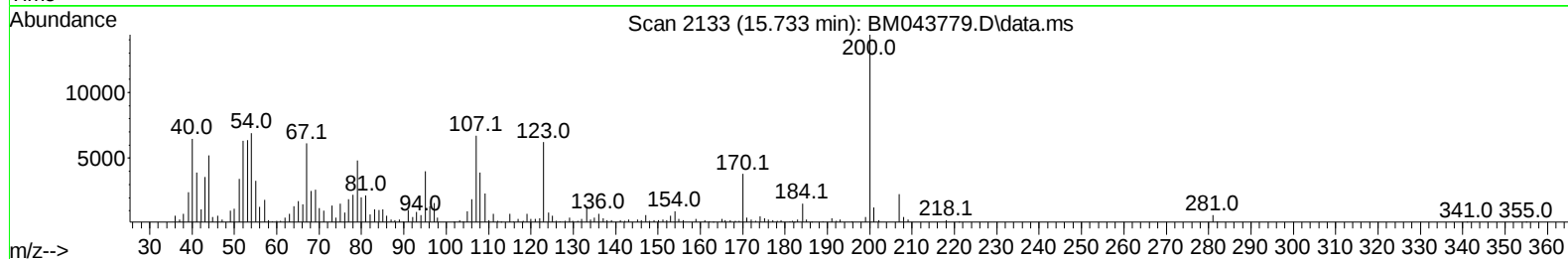
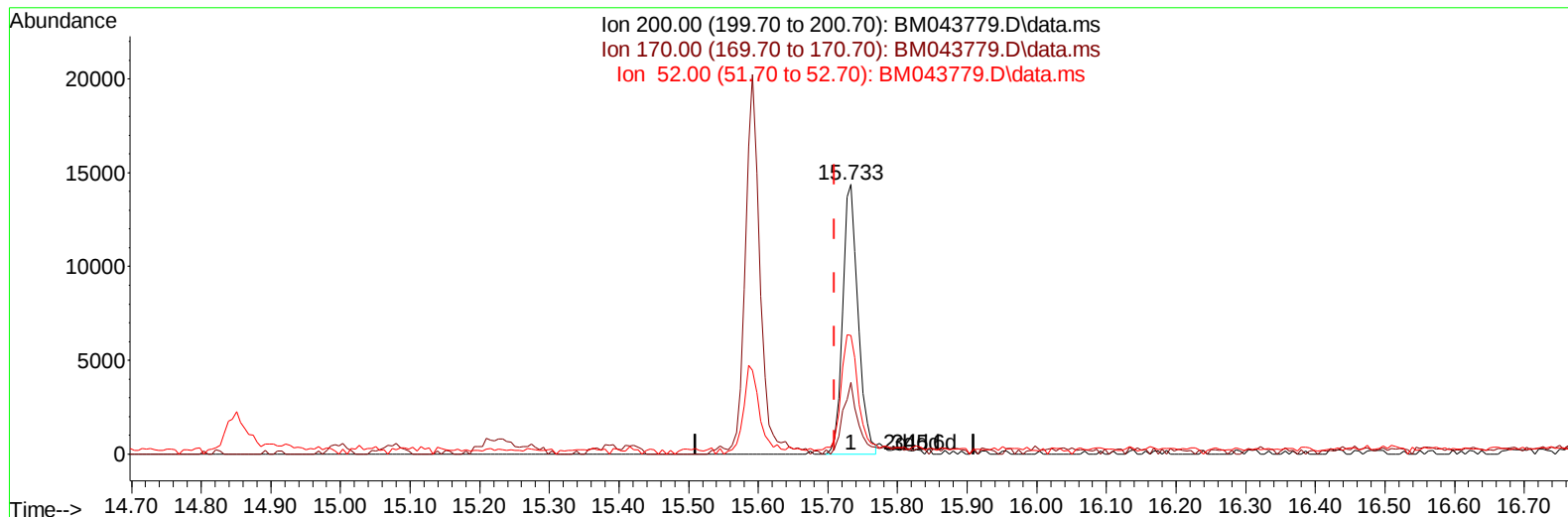
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Data File : BM043779.D
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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BM043779.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.733min (+ 0.023) 2.21 ng/u1

response 22207

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.10	26.44
52.00	58.30	43.91#
0.00	0.00	0.00

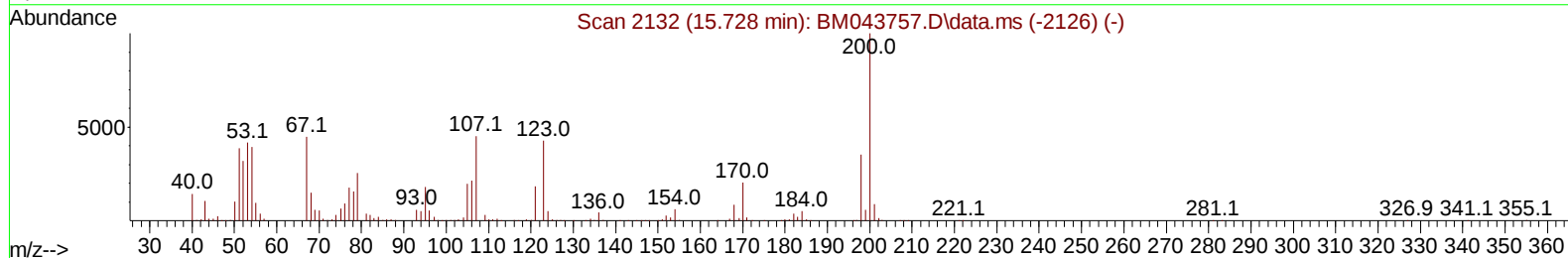
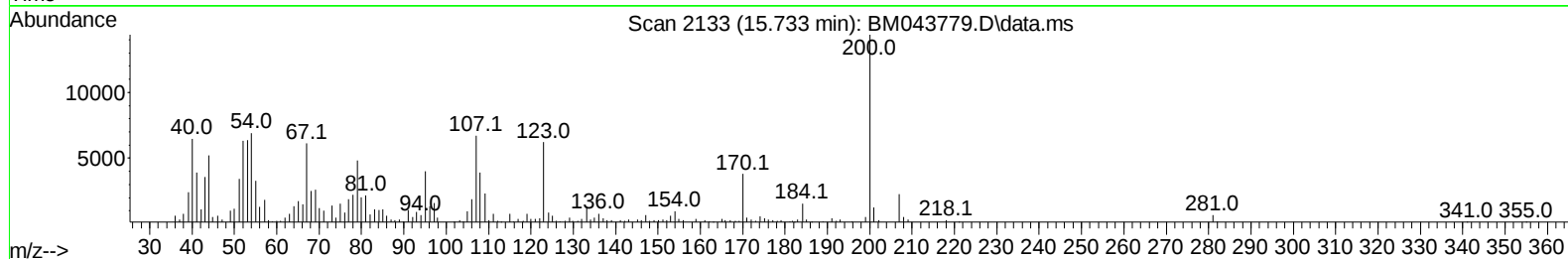
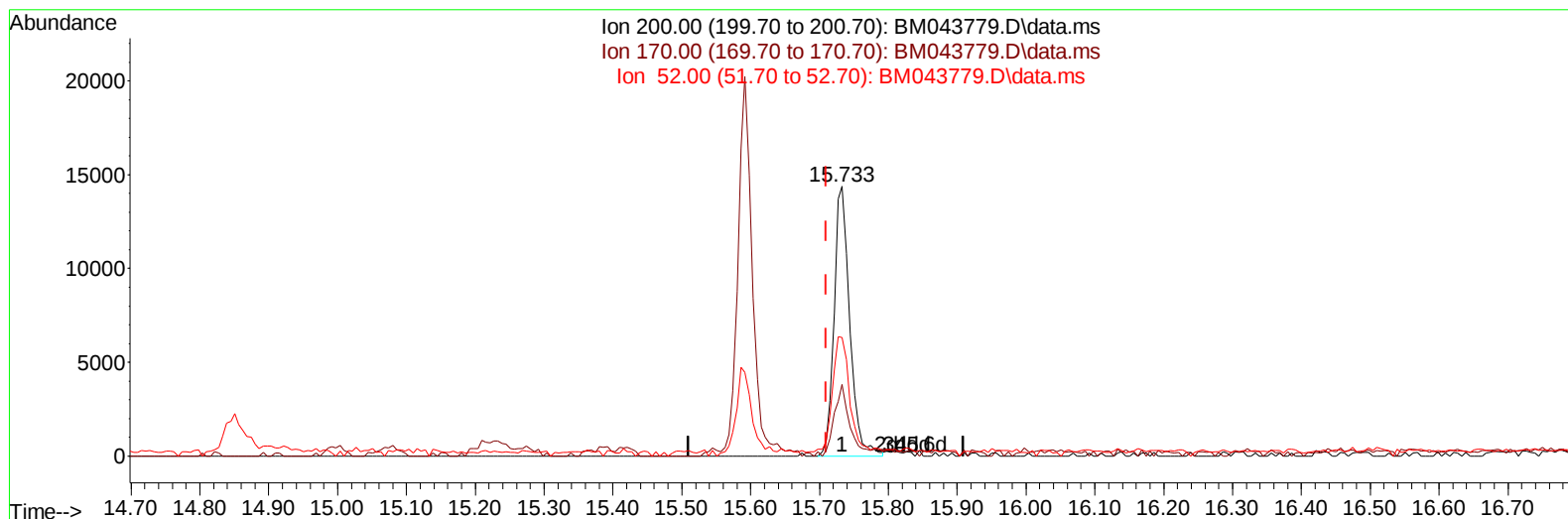
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043779.D
 Acq On : 06 Jan 2024 03:07
 Operator : MA/JU
 Sample : 05952-03MEDL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 01/08/2024



TIC: BM043779.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.733min (+ 0.023) 2.25 ng/ul m

response 22678

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.10	26.44
52.00	58.30	43.91#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/08/2024
 Supervised By :mohammad ahmed 01/08/2024

Quant Time: Jan 06 03:46:02 2024
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.951	152		332869	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136		1310529	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164		736031	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188		1521504	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240		1360835	20.000	ng/ul	0.00
88) Perylene-d12	23.950	264		1582628	20.000	ng/ul	0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.363	96		6656	0.650	ng/uL	0.00
4) Pyridine-d5	3.816	84		31990m	1.085	ng/ul	0.03
7) Phenol-d5	7.134	99		113485	2.997	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.292	67		75032	3.129	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132		93531	3.235	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113		83989	2.858	ng/ul	0.01
21) Nitrobenzene-d5	9.133	128		41564	3.278	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143		44469	3.394	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165		78183	3.381	ng/ul	0.02
31) 4-Chloroaniline-d4	10.927	131		88922	2.613	ng/ul	0.02
46) Dimethylphthalate-d6	14.021	166		253785	3.830	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160		298533	3.798	ng/ul	0.00
54) 4-Nitrophenol-d4	14.845	143		35595m	2.838	ng/ul	0.05
60) Fluorene-d10	15.592	176		228342	4.168	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200		22678m	2.254	ng/ul	0.02
73) Anthracene-d10	17.445	188		351688	4.122	ng/ul	0.00
81) Pyrene-d10	19.739	212		400036	3.723	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.791	264		404543	4.142	ng/ul	0.02
Target Compounds							
						Qvalue	
71) Pentachlorophenol	16.998	266		551832	44.200	ng/ul	99

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

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