

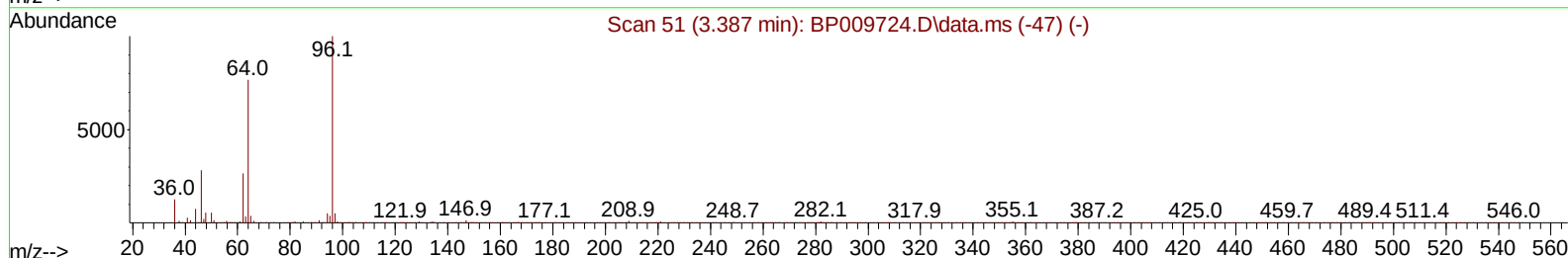
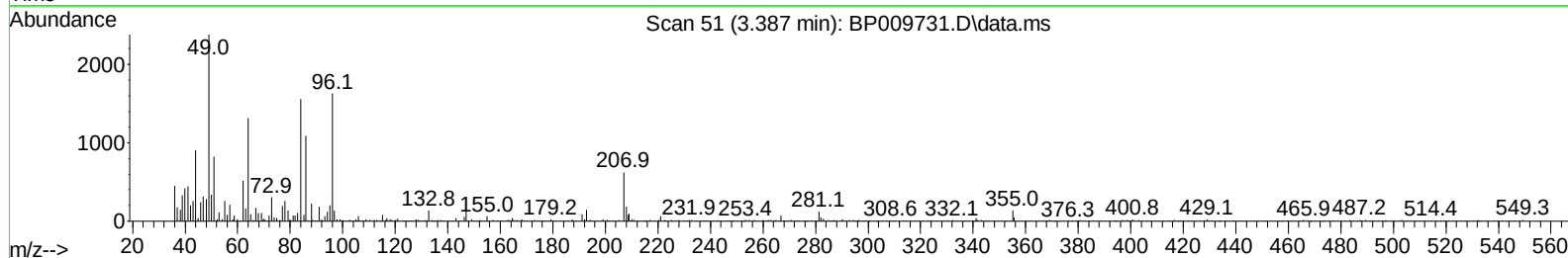
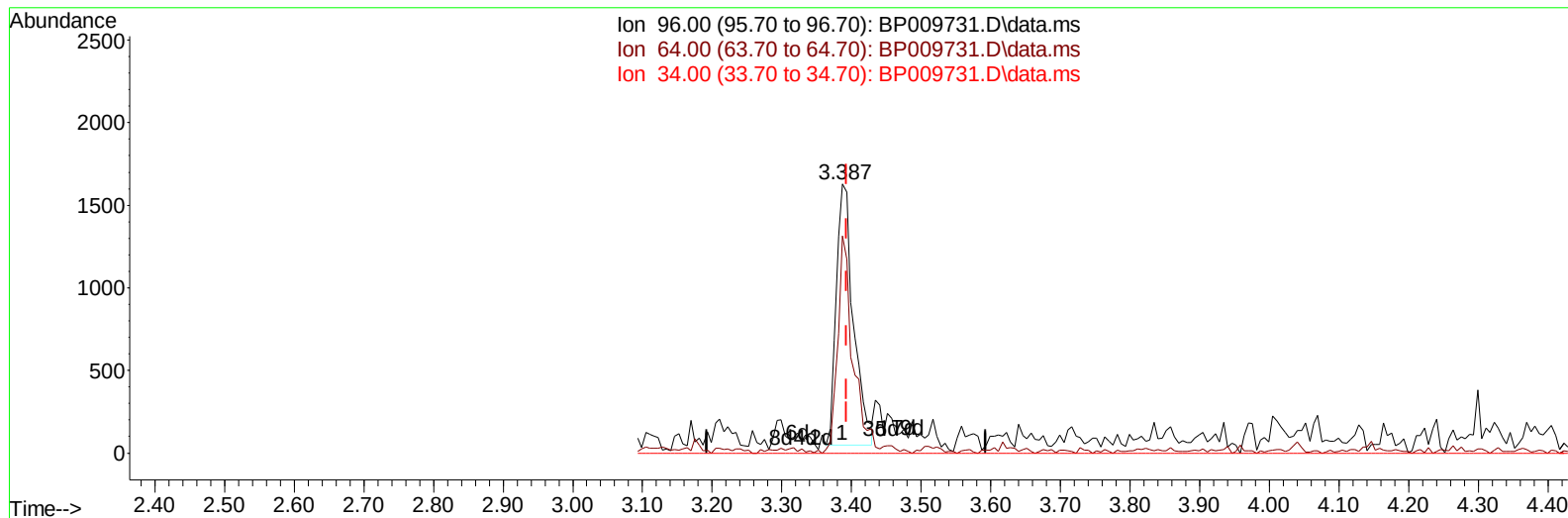
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009731.D
 Acq On : 08 Apr 2022 23:17
 Operator : CG/JU
 Sample : N2182-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R14

Manual IntegrationsAPPROVED

Quant Time: Apr 09 01:16:31 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: BP009731.D\data.ms

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

3.387min (-0.006) 2.80 ng/uL

response 2690

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.65#
34.00	0.00	0.00
0.00	0.00	0.00

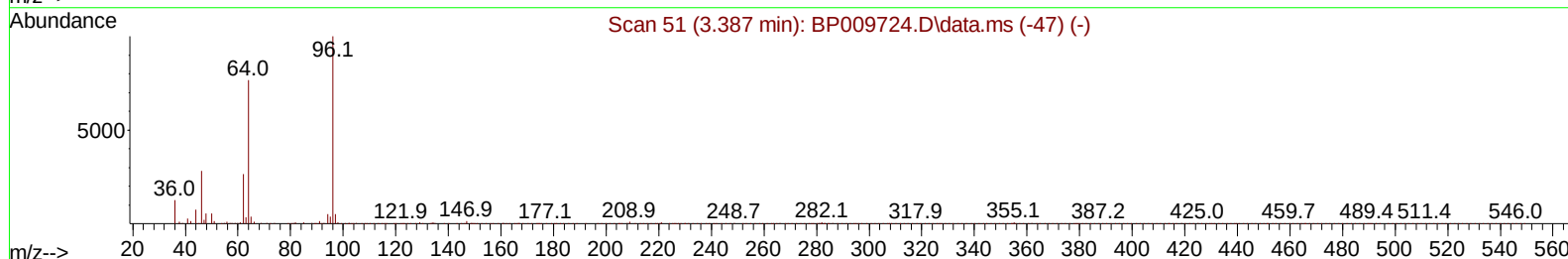
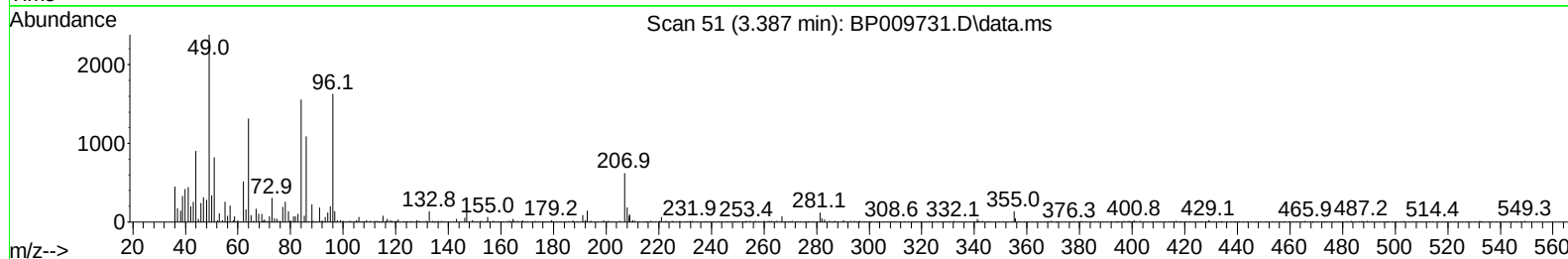
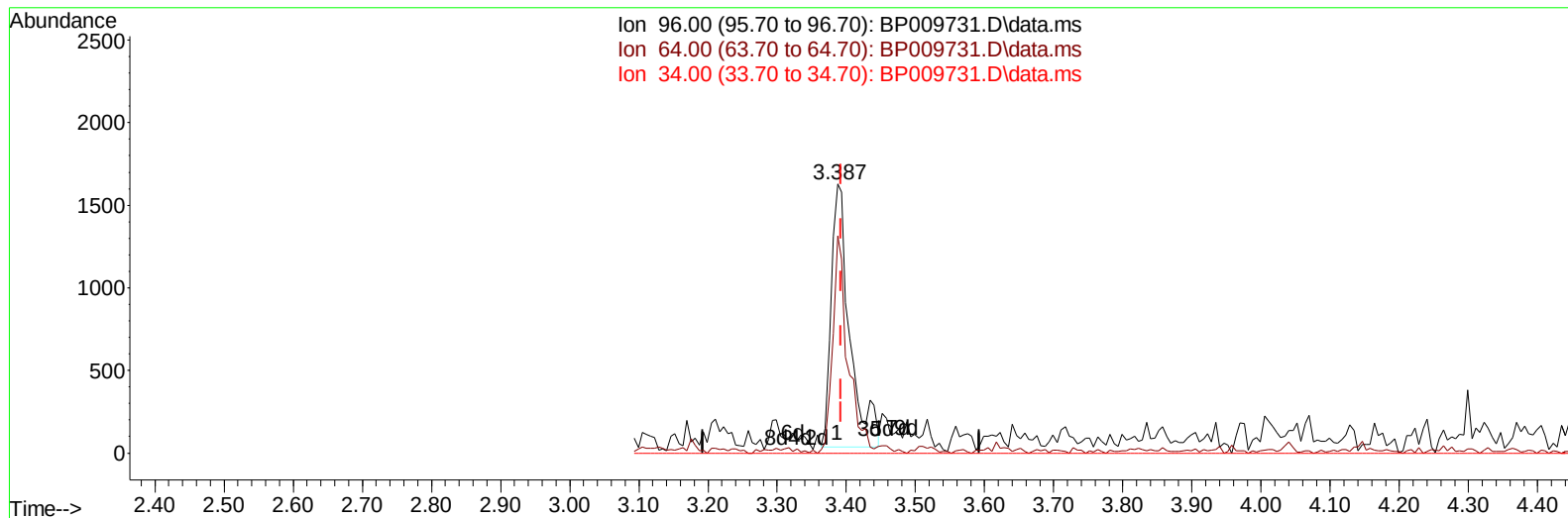
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TIC: BP009731.D\data.ms

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

3.387min (-0.006) 3.07 ng/uL m

response 2949

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.65#
34.00	0.00	0.00
0.00	0.00	0.00

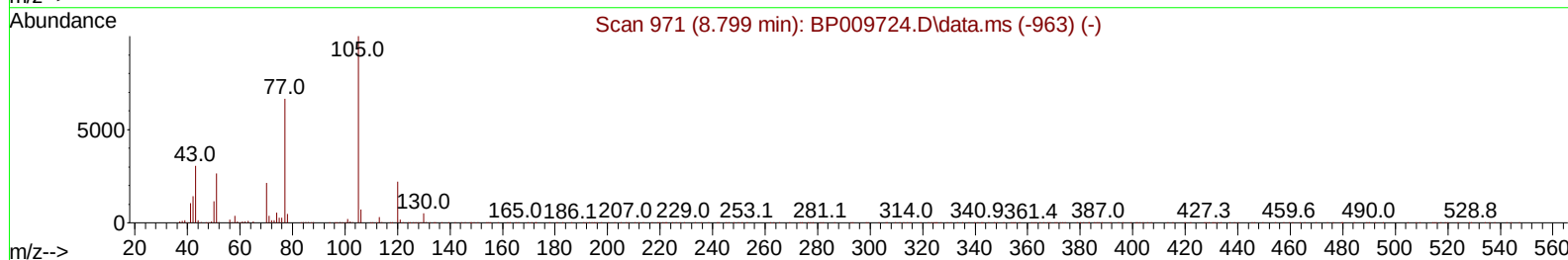
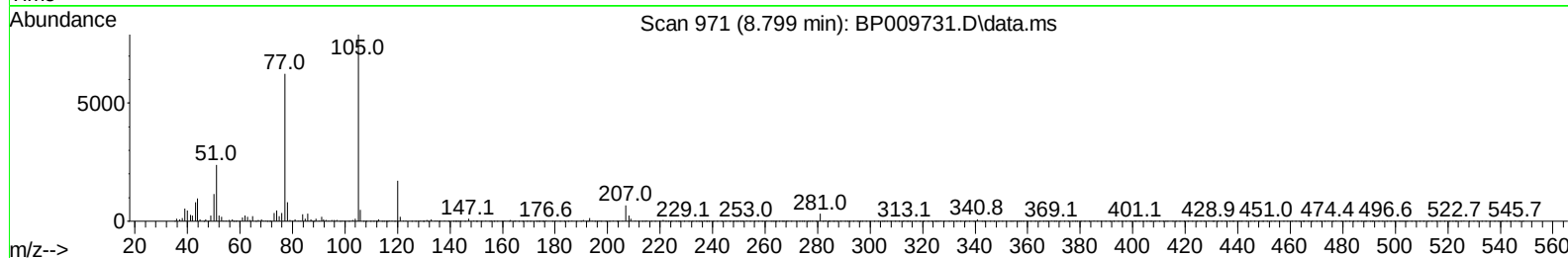
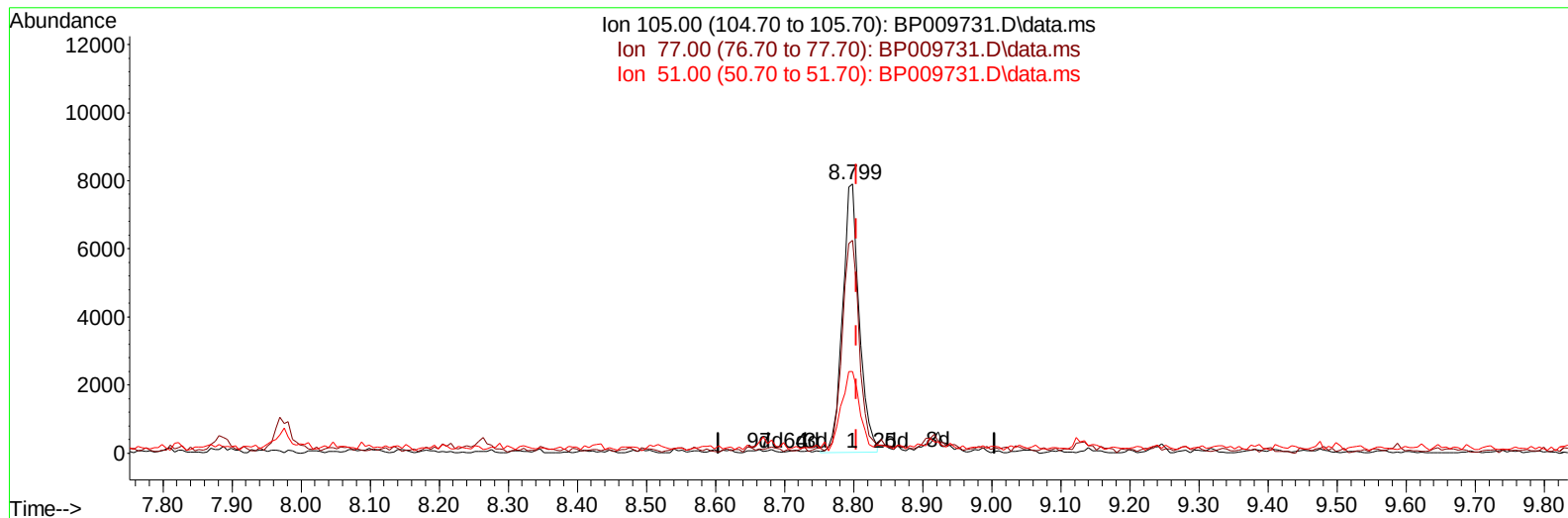
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009731.D
 Acq On : 08 Apr 2022 23:17
 Operator : CG/JU
 Sample : N2182-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP009731.D\data.ms

(16) Acetophenone

8.799min (-0.006) 2.68 ng/u1

response 13257

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	79.00
51.00	17.30	30.27#
0.00	0.00	0.00

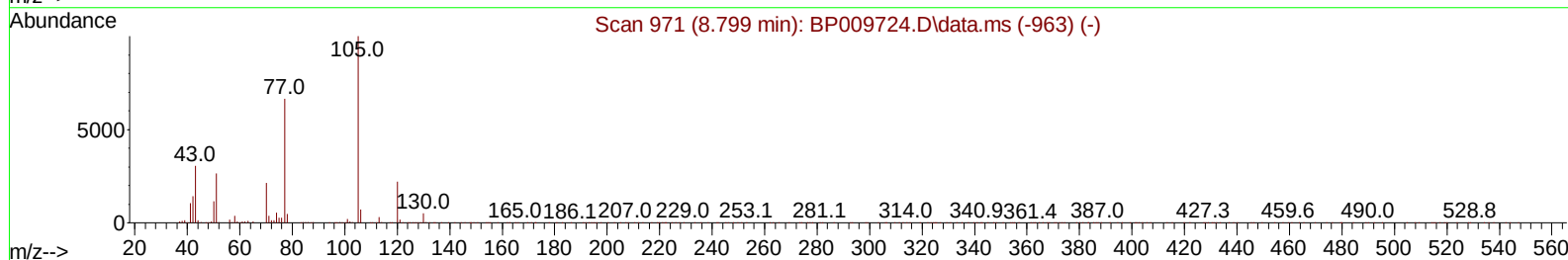
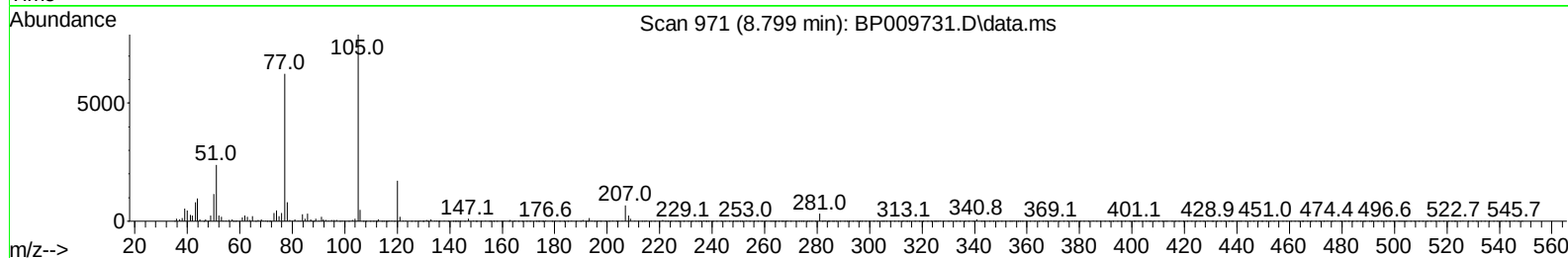
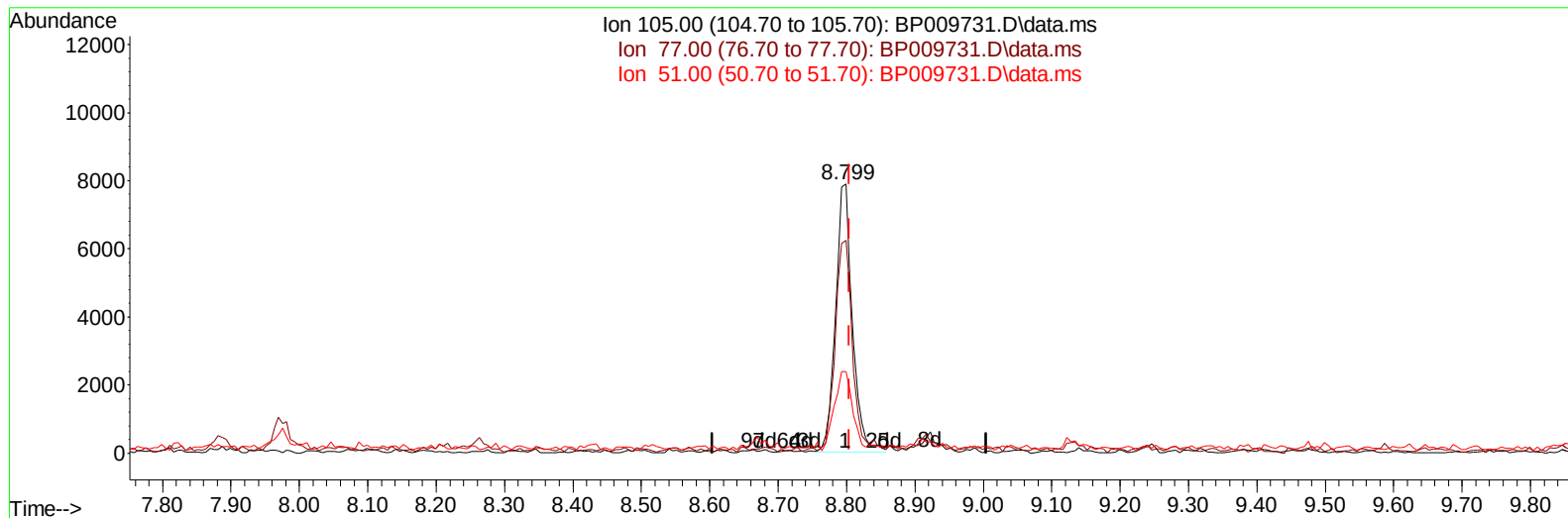
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
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TIC: BP009731.D\data.ms

(16) Acetophenone

8.799min (-0.006) 2.70 ng/ul m

response 13370

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	79.00
51.00	17.30	30.27#
0.00	0.00	0.00

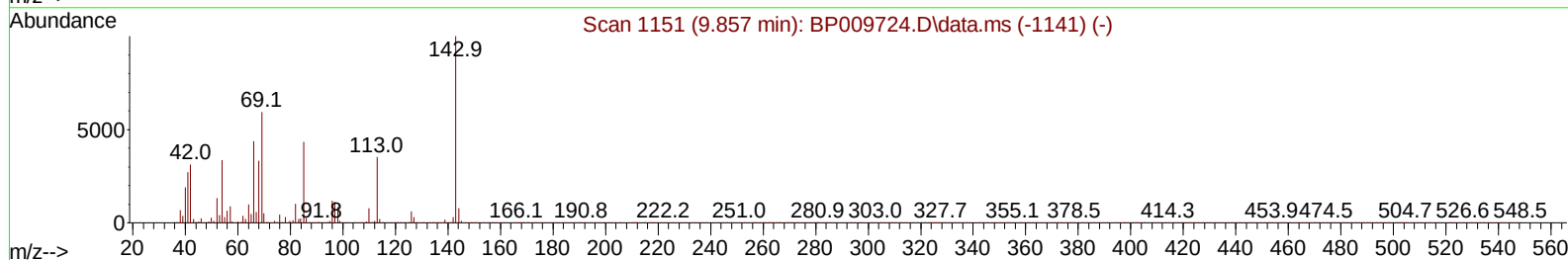
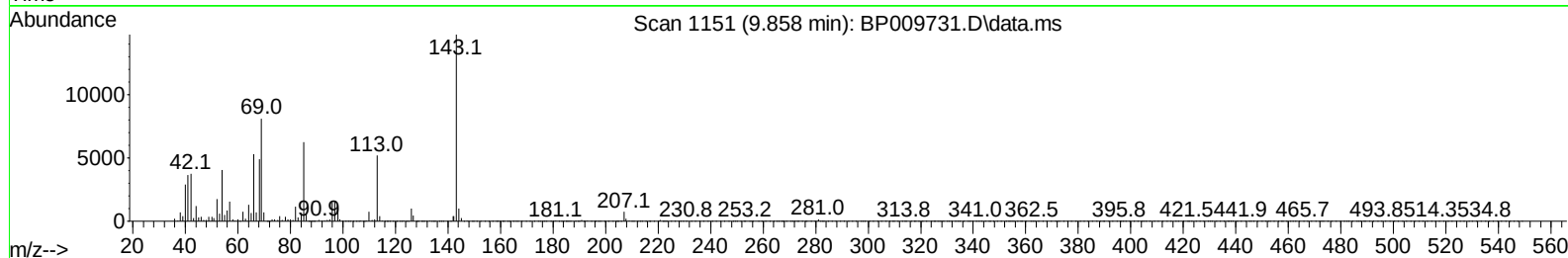
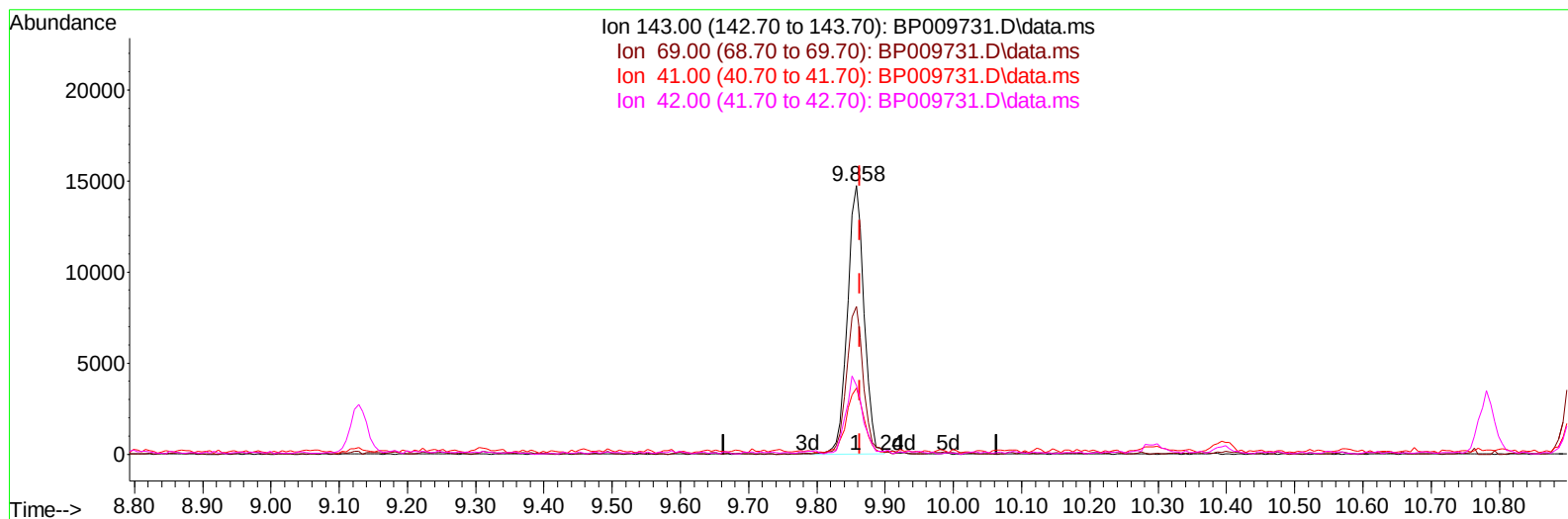
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009731.D
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 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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TIC: BP009731.D\data.ms

(24) 2-Nitrophenol-d4 (S)

9.858min (-0.006) 16.46 ng/ul

response 25001

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	47.10	55.05
41.00	16.70	24.63#
42.00	20.70	25.59#

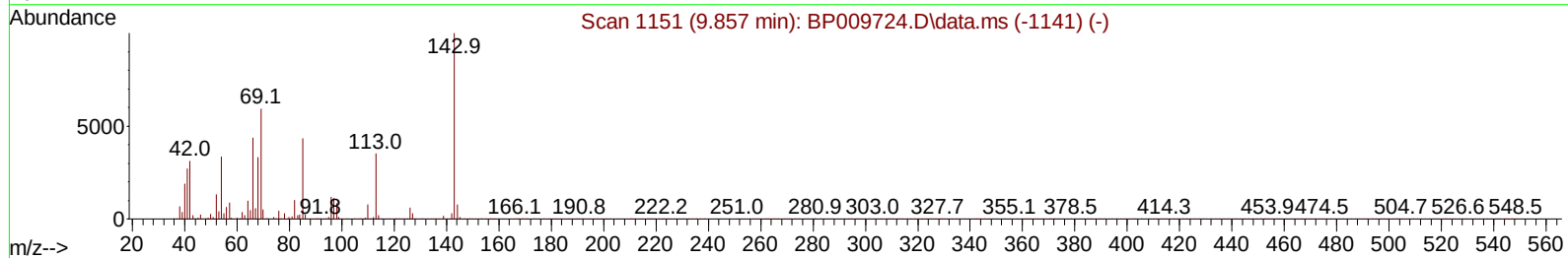
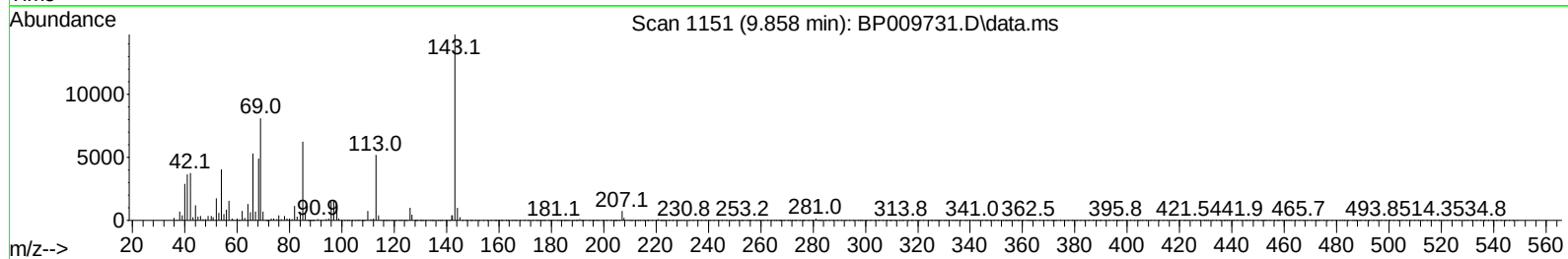
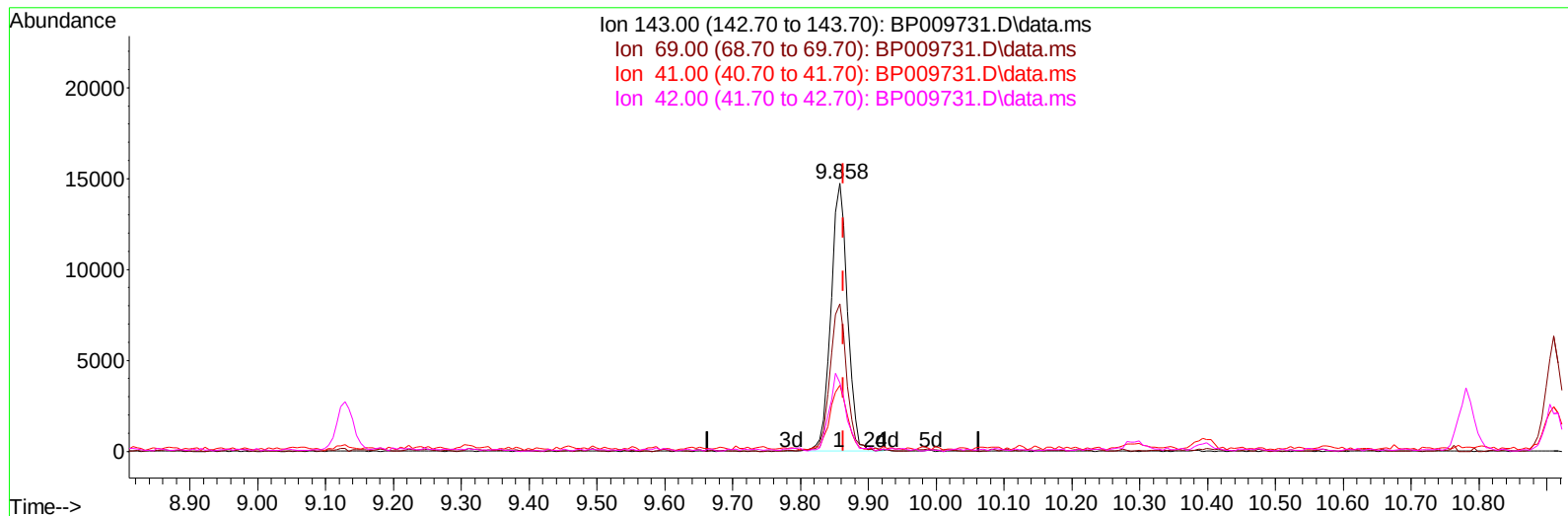
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009731.D
 Acq On : 08 Apr 2022 23:17
 Operator : CG/JU
 Sample : N2182-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R14

Manual IntegrationsAPPROVED

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TIC: BP009731.D\data.ms

(24) 2-Nitrophenol-d4 (S)

9.858min (-0.006) 16.42 ng/ul m

response 24947

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	47.10	55.05
41.00	16.70	24.63#
42.00	20.70	25.59#

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 Operator : CG/JU
 Sample : N2182-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.975	152		42845	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136		192168	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164		144833	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188		341924	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240		411116	20.000	ng/ul	#-0.01
88) Perylene-d12	23.910	264		482153	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.387	96		2949m	3.072	ng/uL	0.00
4) Pyridine-d5	3.817	84		8313	3.141	ng/ul	0.00
7) Phenol-d5	7.122	99		57494	15.285	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67		37704	16.837	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132		45018	15.866	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113		38212	12.169	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128		23596	17.027	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.858	143		24947m	16.420	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165		50956	16.028	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131		57119	12.614	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166		212077	18.641	ng/ul	-0.01
49) Acenaphthylene-d8	14.299	160		221205	16.412	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143		35156	17.039	ng/ul	-0.01
60) Fluorene-d10	15.593	176		183674	19.191	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.704	200		24735	12.211	ng/ul	0.00
73) Anthracene-d10	17.451	188		318645	19.464	ng/ul	-0.01
81) Pyrene-d10	19.681	212		450769	20.076	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264		525960	20.947	ng/ul	-0.01
Target Compounds							
16) Acetophenone	8.799	105		13370m	2.702	ng/ul	Qvalue

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

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