

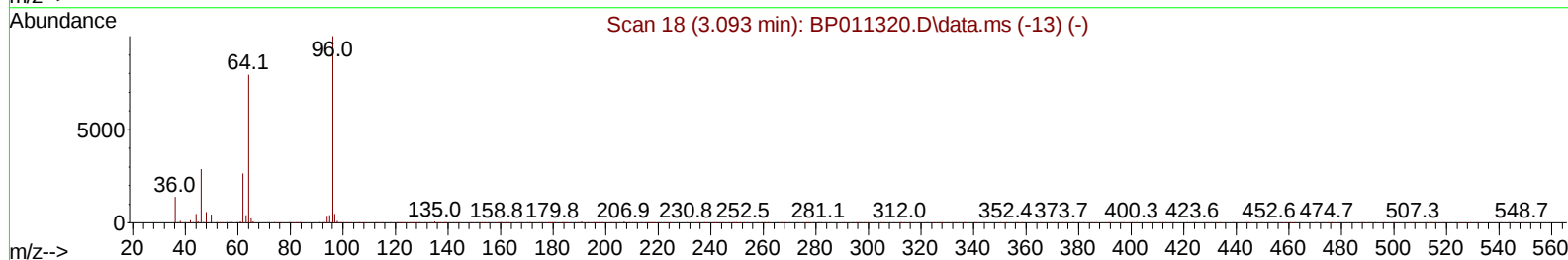
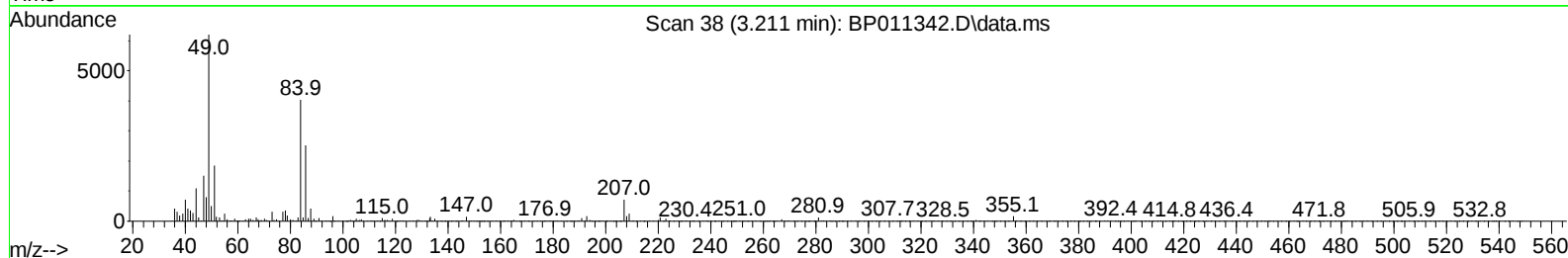
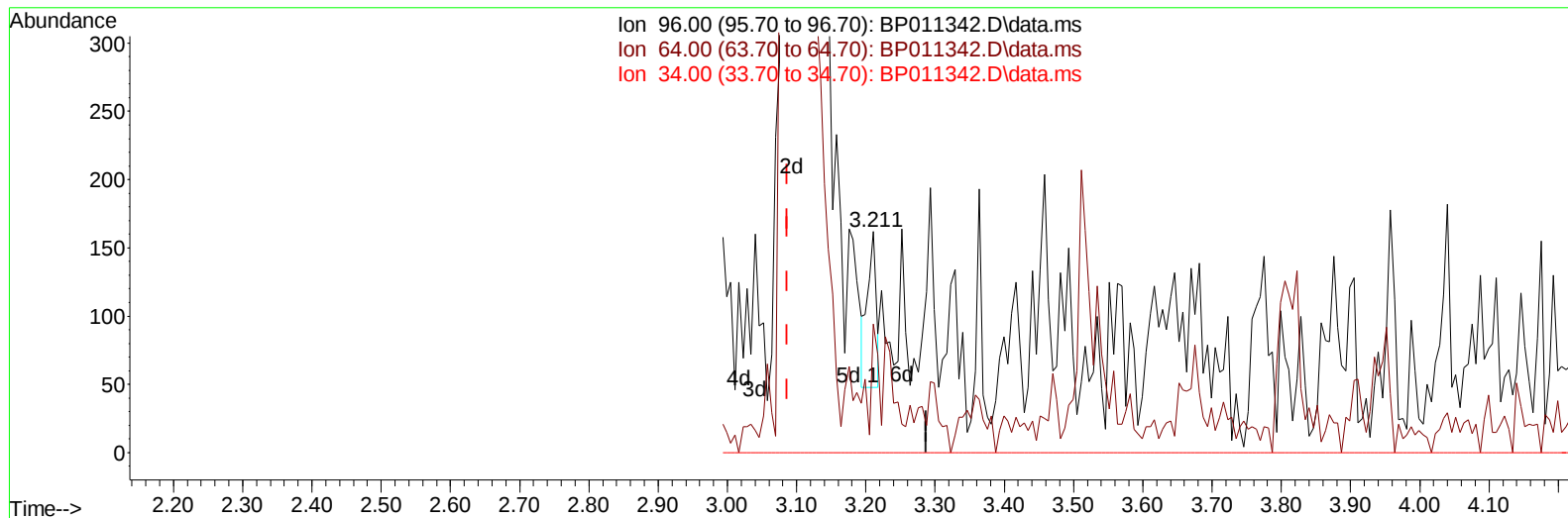
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011342.D
Acq On : 09 Aug 2022 22:29
Operator : CG/JU
Sample : N3900-04
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0C94

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:17:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011342.D\data.ms

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

3.211min (+ 0.124) 0.03 ng/uL

response	101
Ion	Exp% Act%
96.00	100.00 100.00
64.00	61.00 54.65
34.00	0.00 0.00
0.00	0.00 0.00

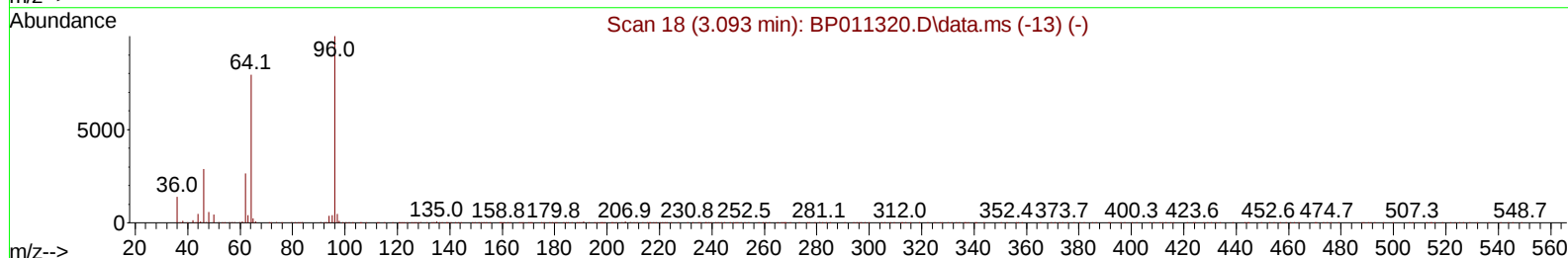
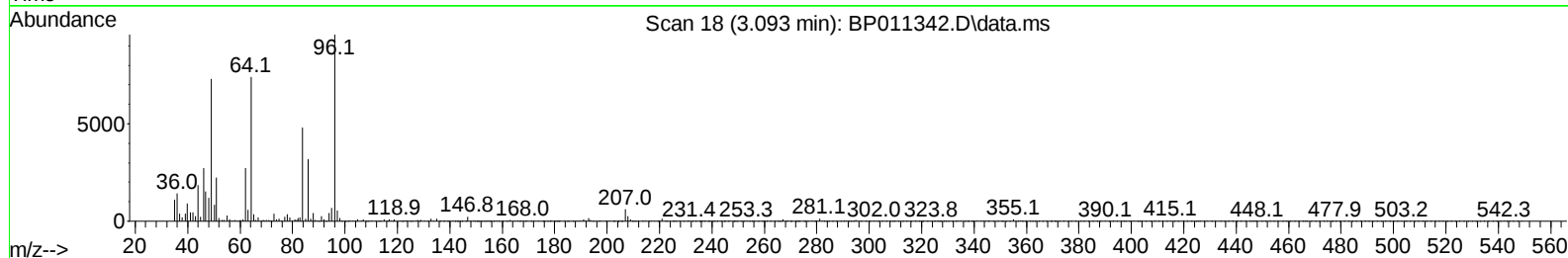
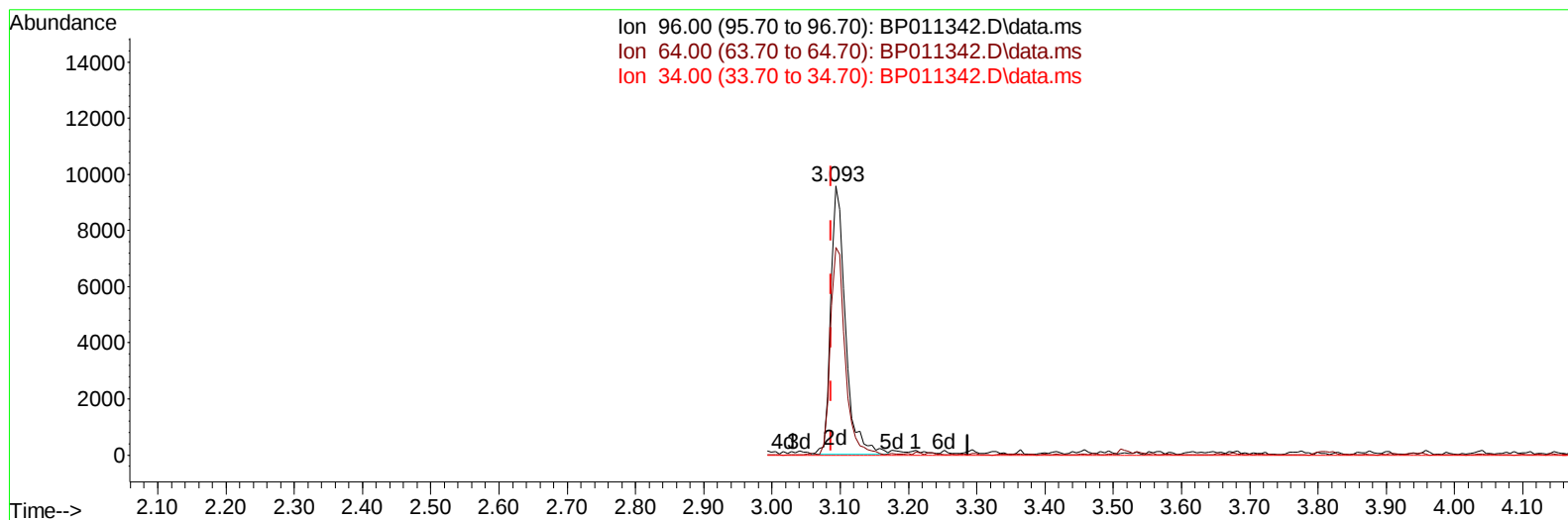
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(3) 1,4-Dioxane-d8 (S)

3.093min (+ 0.006) 3.86 ng/uL m

response 14282

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	77.14#
34.00	0.00	0.00
0.00	0.00	0.00

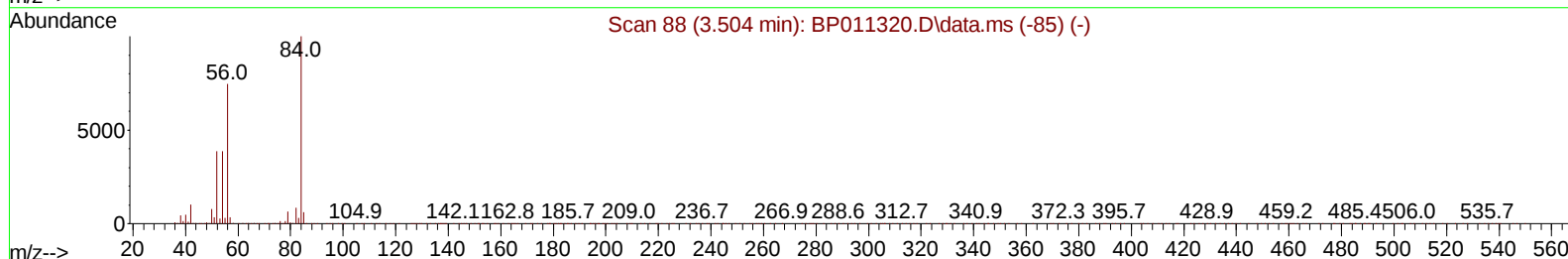
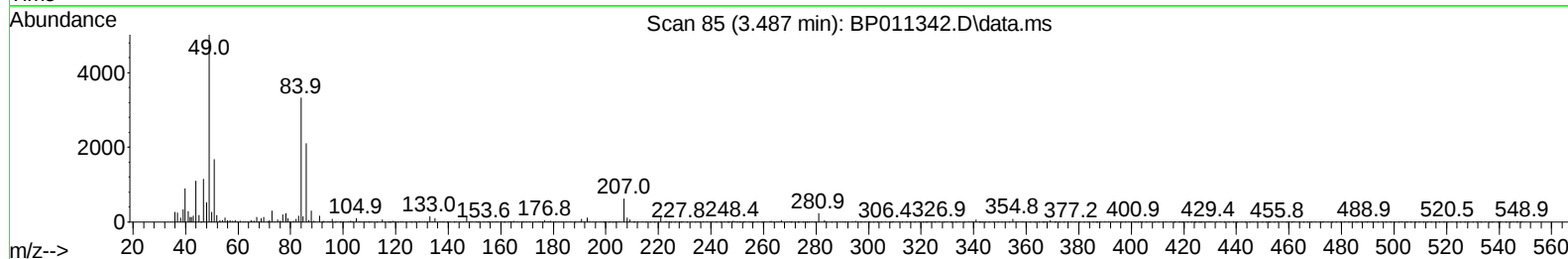
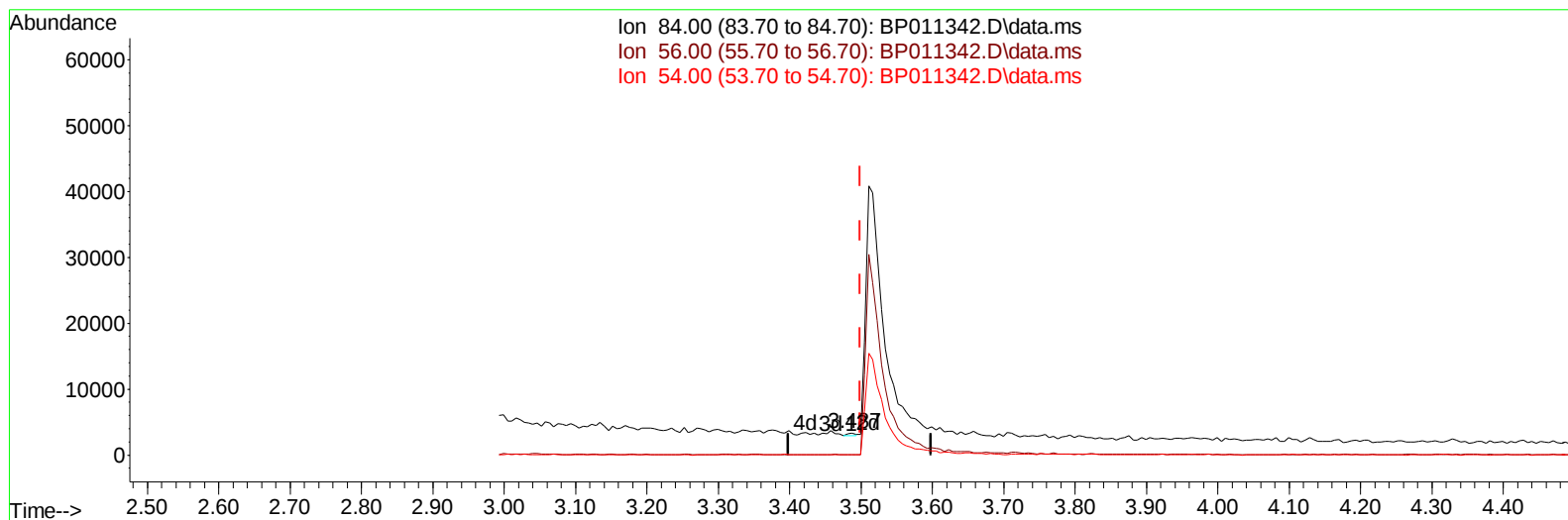
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TIC: BP011342.D\data.ms

(4) Pyridine-d5 (S)

3.487min (-0.012) 0.03 ng/u1

response 286

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	1.44#
54.00	33.20	1.56#
0.00	0.00	0.00

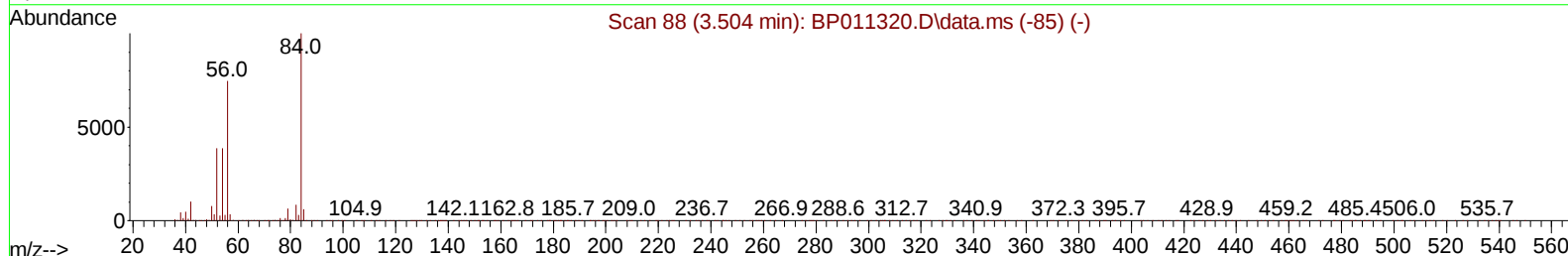
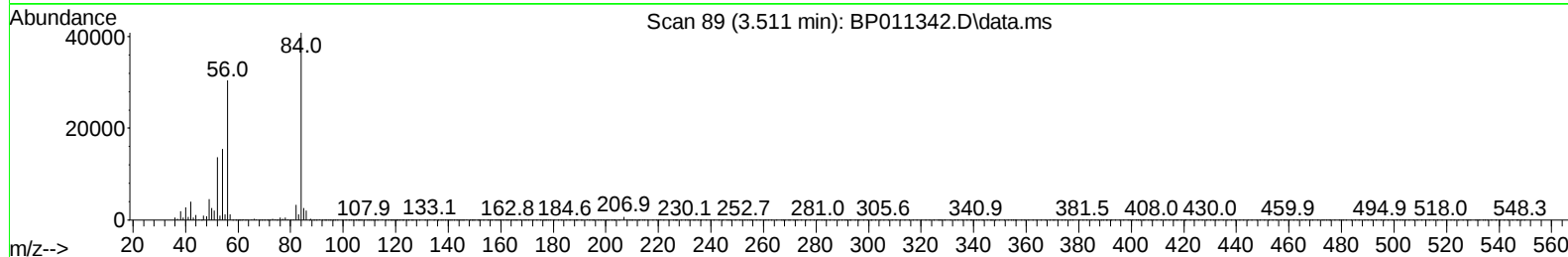
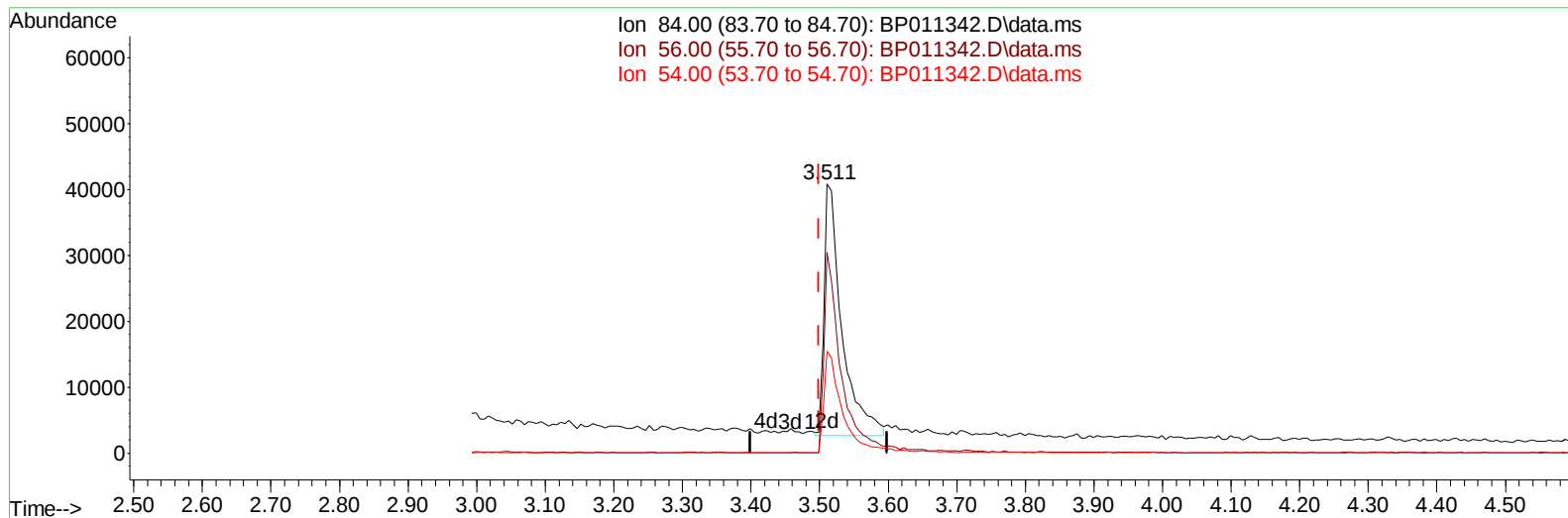
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TIC: BP011342.D\data.ms

(4) Pyridine-d5 (S)

3.511min (+ 0.012) 7.33 ng/u1 m

response 68532

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	74.55#
54.00	33.20	37.89
0.00	0.00	0.00

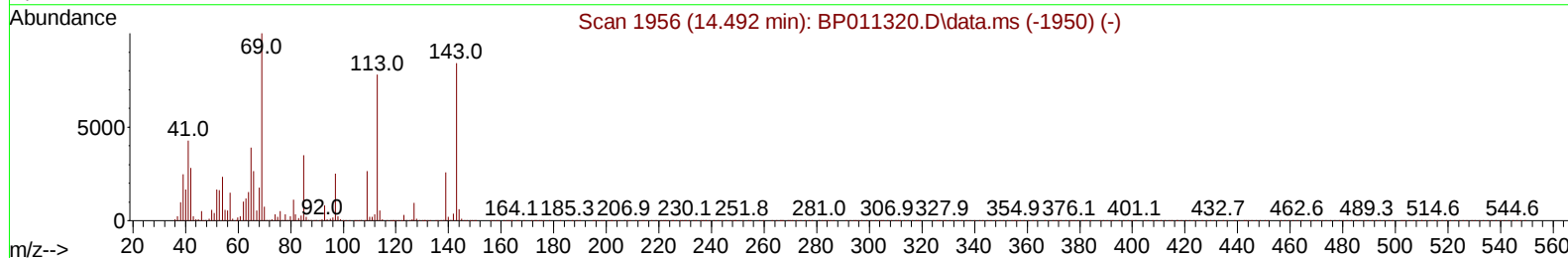
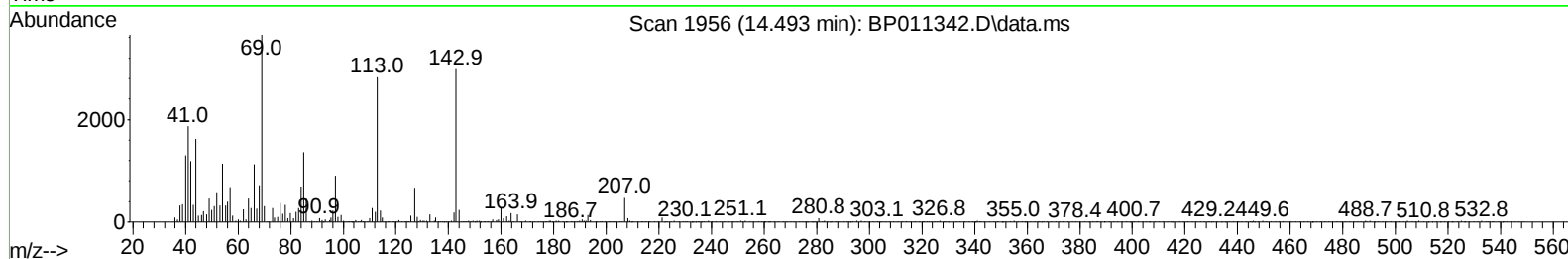
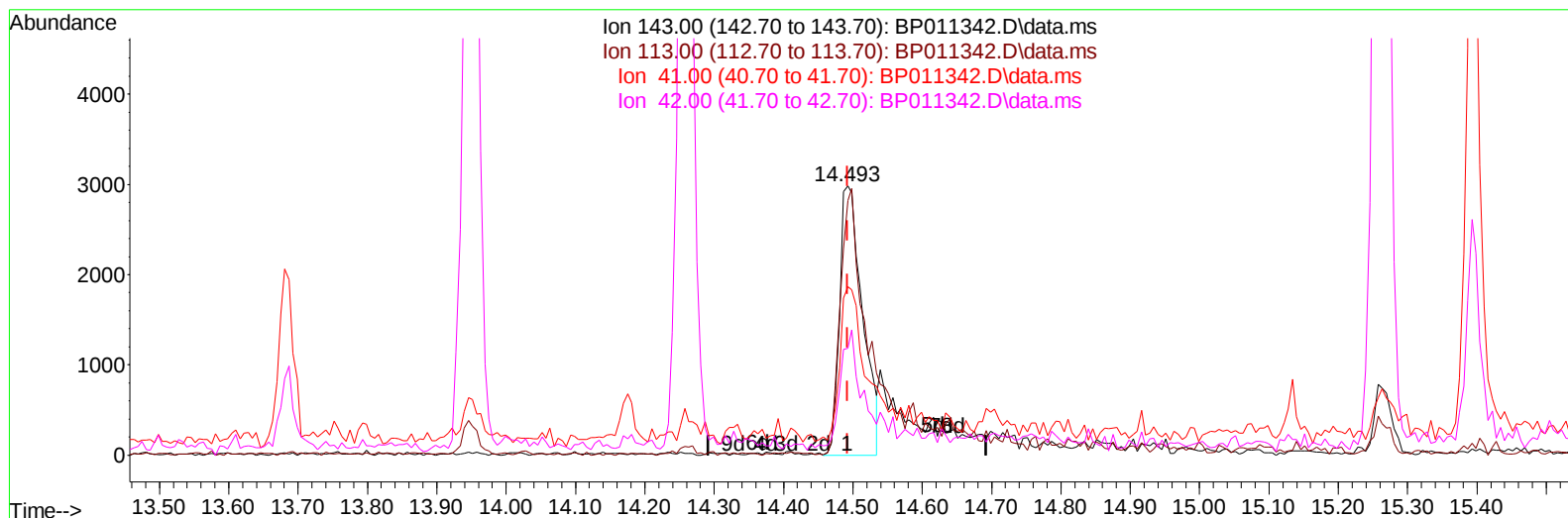
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TIC: BP011342.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.000) 1.96 ng/u1

response 6800

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.73#
41.00	54.40	62.66
42.00	35.50	39.72

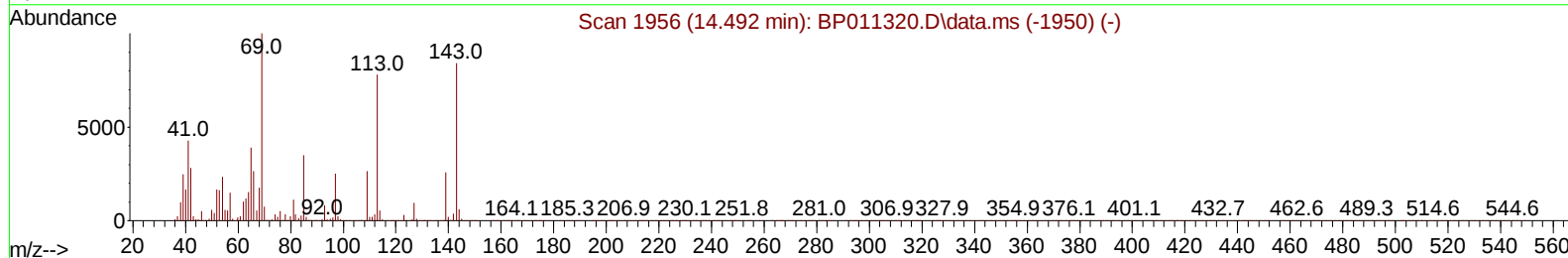
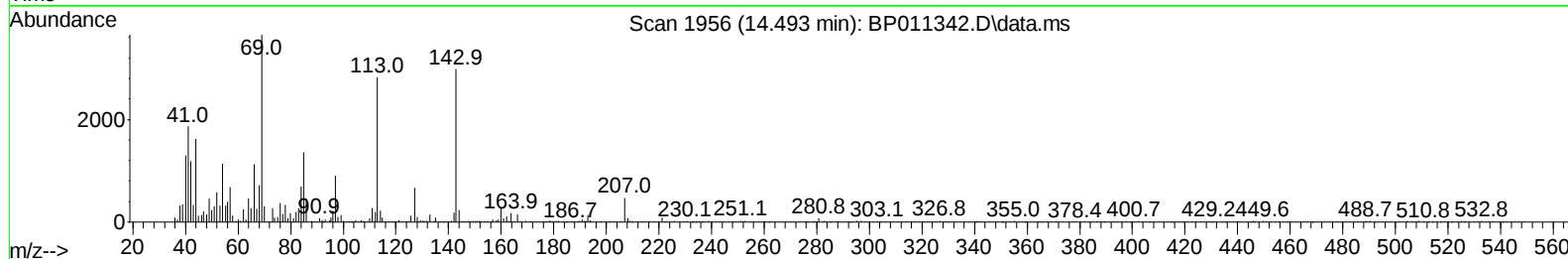
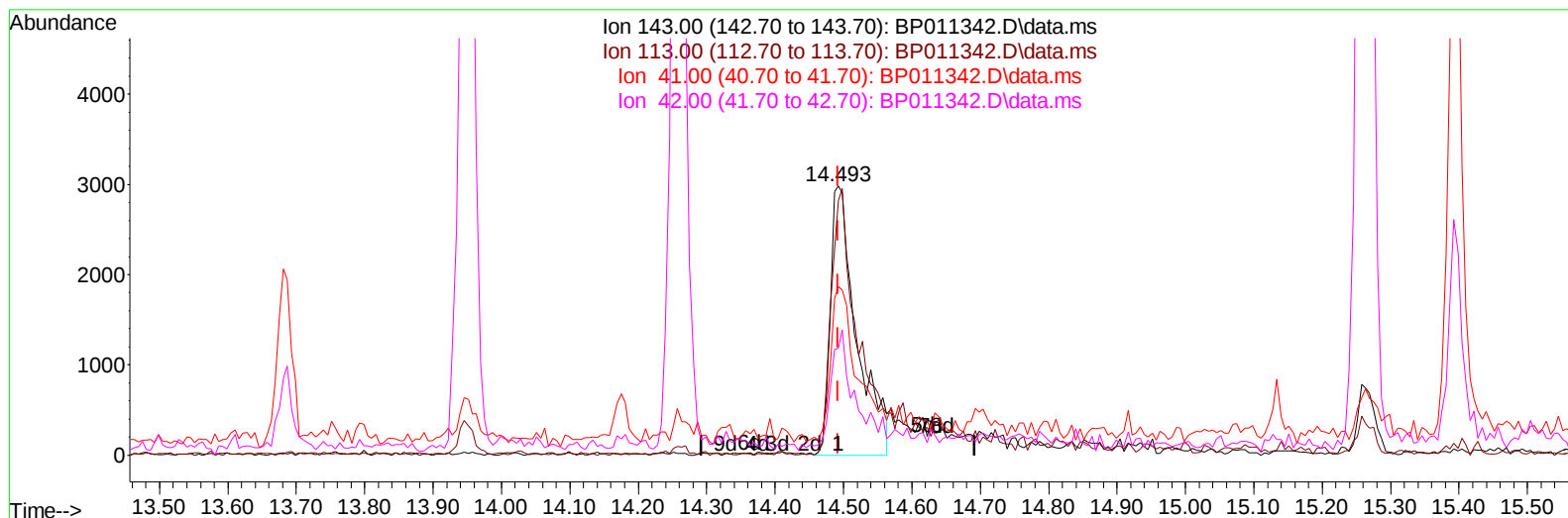
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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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TIC: BP011342.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.000) 2.29 ng/ul m

response 7946

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.73#
41.00	54.40	62.66
42.00	35.50	39.72

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	113568	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	475512	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	274354	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	543001	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	509387	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	515023	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	14282m	3.855	ng/uL	0.00
4) Pyridine-d5	3.511	84	68532m	7.334	ng/ul	0.01
7) Phenol-d5	6.775	99	47972	4.773	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	190851	26.781	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	162562	20.614	ng/ul	0.00
15) 4-Methylphenol-d8	8.311	113	94640	11.993	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	100777	31.140	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	91617	30.921	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	151680	25.350	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	239366	22.882	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	615161	32.316	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	707739	30.886	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	7946m	2.285	ng/ul	0.00
60) Fluorene-d10	15.263	176	522136	32.073	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	67512	26.767	ng/ul	0.00
73) Anthracene-d10	17.134	188	864843	35.435	ng/ul	0.00
81) Pyrene-d10	19.392	212	993338	35.395	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	925230	35.623	ng/ul	-0.01

Target Compounds Qvalue

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

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