

Legends for Figures

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Figure S2: ^{13}C NMR spectrum of friedelan-3-one (**1**) recorded in CDCl_3 .

Figure S3: MS spectrum of friedelan-3-one (**1**).

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Figure S5: ^1H NMR spectrum of 3α -Hydroxyfriedel-2-one (**2**) recorded in CDCl_3 .

Figure S6: ^{13}C NMR spectrum of 3α -Hydroxyfriedel-2-one (**2**) recorded in CDCl_3 .

Figure S7: MS spectrum of 3α -Hydroxyfriedel-2-one (**2**).

Figure S8: FT-IR spectrum of 3α -Hydroxyfriedel-2-one (**2**).

Figure S9: ^1H NMR spectrum of 3-Hydroxyfriedel-3-en-2-one (**3**) recorded in CDCl_3 .

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Figure S13: ^1H NMR spectrum of Lupeol (**4**) recorded in CDCl_3 .

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Figure S25. ^1H NMR spectrum of Oleanolic acid acetate (**7**) recorded in CDCl_3 .

Figure S26: ^{13}C NMR spectrum of Oleanolic acid acetate (**7**) recorded in CDCl_3

Figure S27: MS spectrum of Oleanolic acid acetate (**7**).

Figure S28: FT-IR spectrum of Oleanolic acid acetate (**7**).

Figure S29: ^1H NMR spectrum of Tetradecyl (E)-ferulate (**8**) recorded in CDCl_3 .

Figure S30: ^{13}C NMR spectrum of Tetradecyl (E)-ferulate (**8**) recorded in CDCl_3

Figure S31: MS spectrum of Tetradecyl (E)-ferulate (**8**).

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Figure S33: ^1H NMR spectrum of 9-Octadecenoic acid (**9**) recorded in CDCl_3 .

Figure S34: ^{13}C NMR spectrum of 9-Octadecenoic acid (**9**) recorded in CDCl_3

Figure S35: MS spectrum of 9-Octadecenoic acid (**9**).

Figure S36: FT-IR spectrum of 9-Octadecenoic acid (**9**).

Table S1: ^{13}C NMR (100 MHz) data (δ value) of isolated compounds (1-8) in CHCl_3

Table S2: ^1H NMR (400 MHz) data (δ value) of isolated compounds (1-5 and 7) in CHCl_3

Table S3: ^1H NMR (400 MHz) data (δ value) of isolated compounds (6 and 8) in CHCl

PROTON_07
ZTPA80-95_major_cult

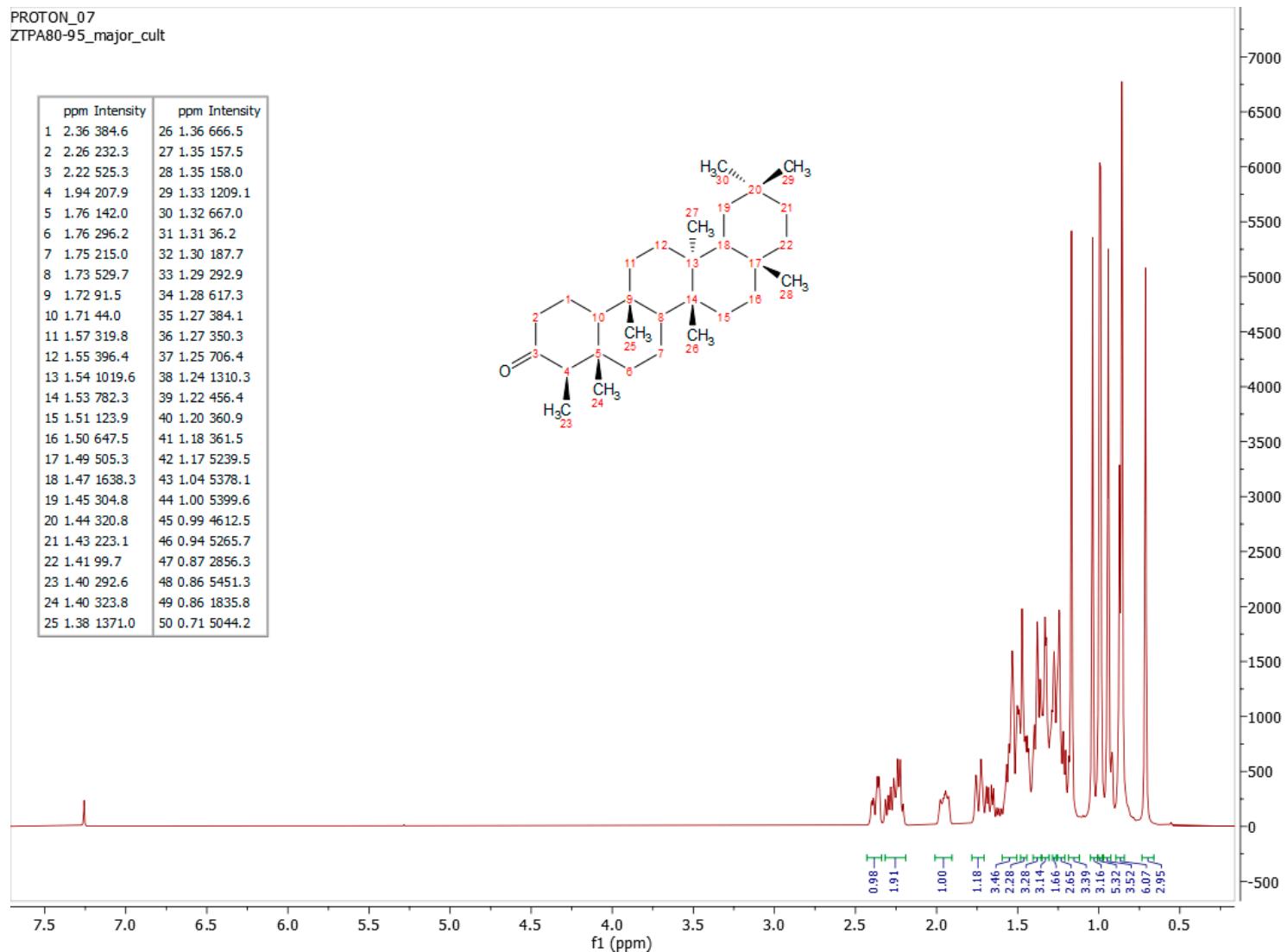


Figure SI: ¹H NMR spectrum of friedelan-3-one (1) recorded in CDCl₃.

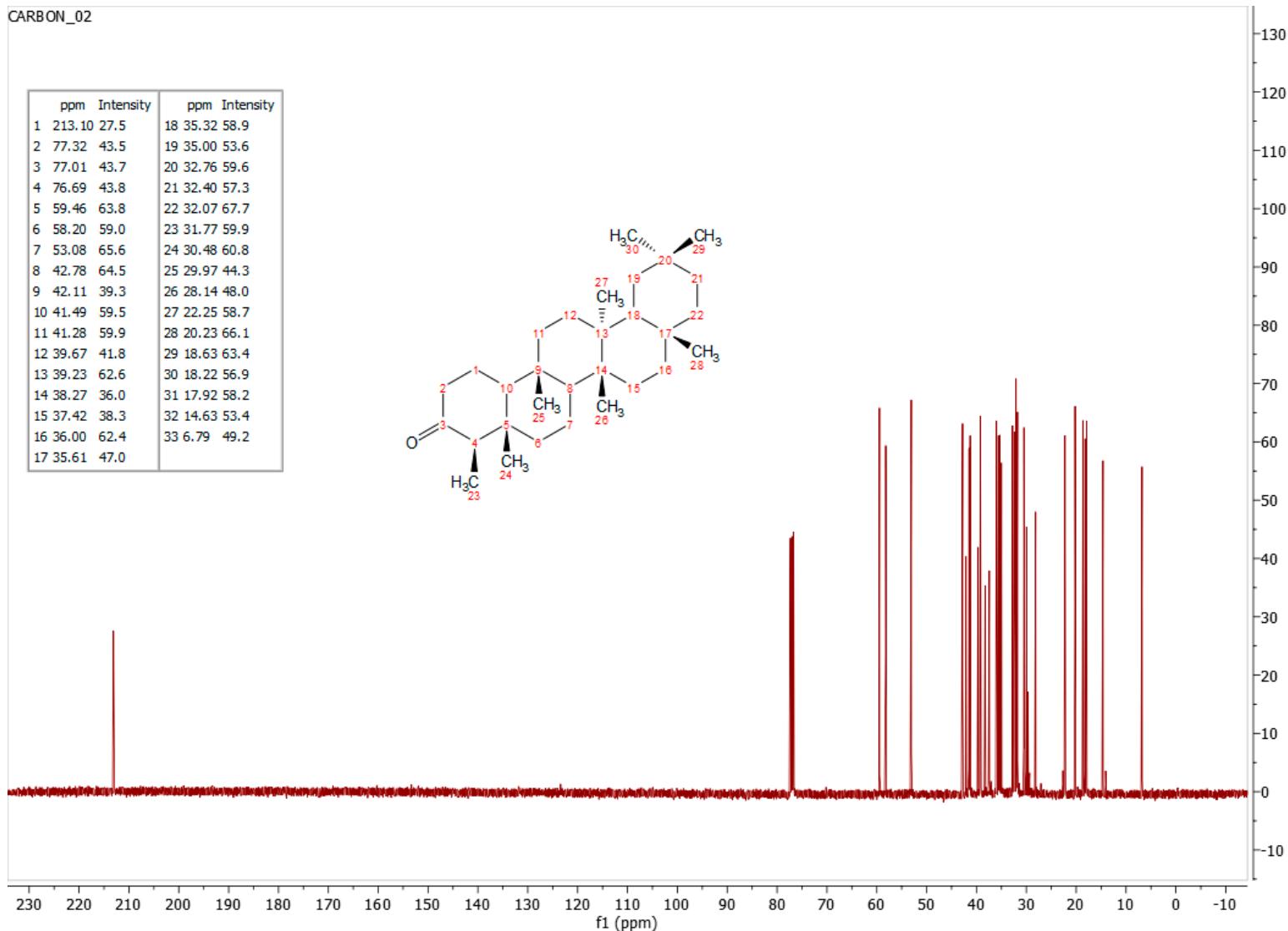


Figure S2: ¹³C NMR spectrum of friedelan-3-one (**1**) recorded in CDCl₃.

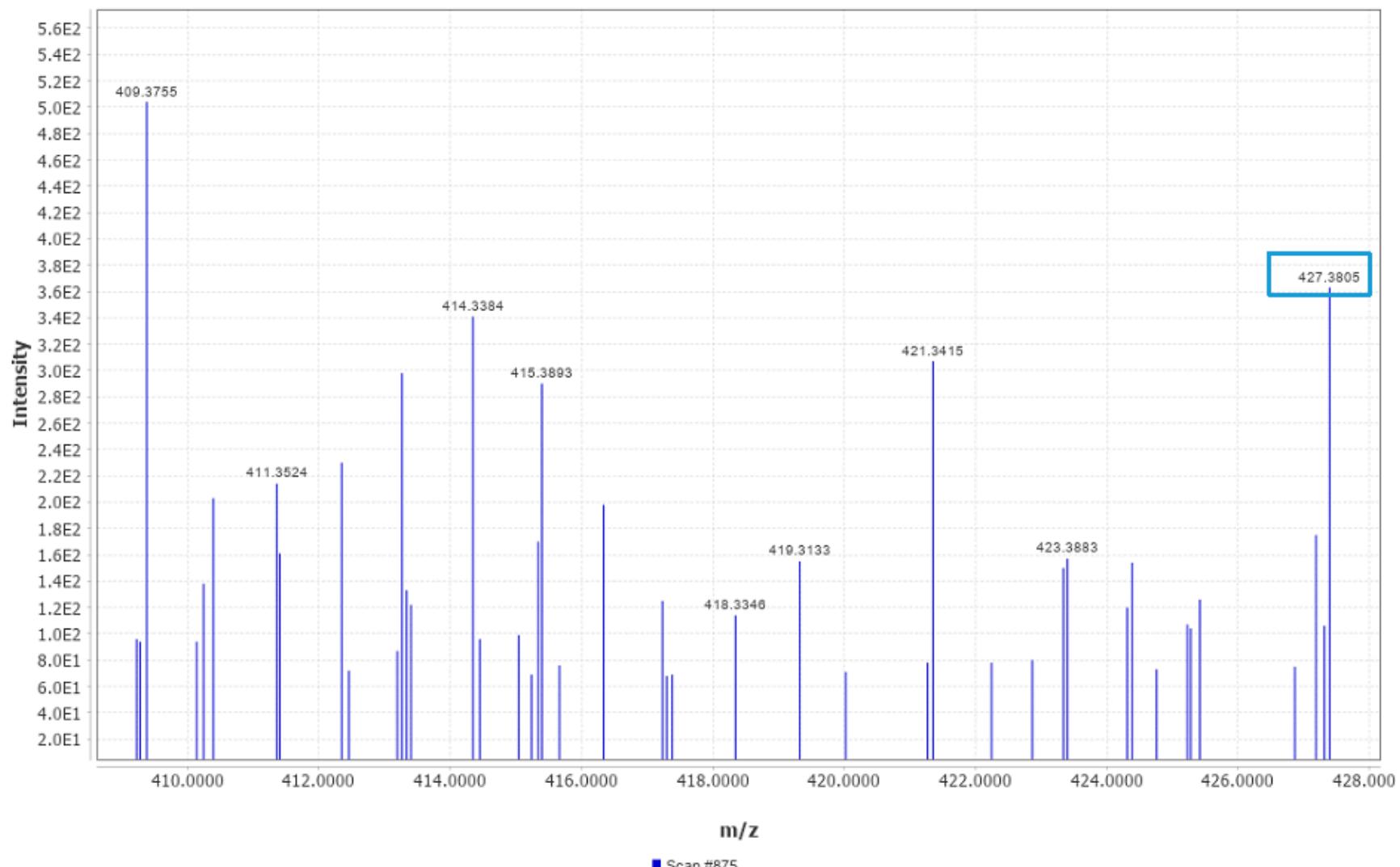


Figure S3: MS spectrum of friedelan-3-one (**1**).

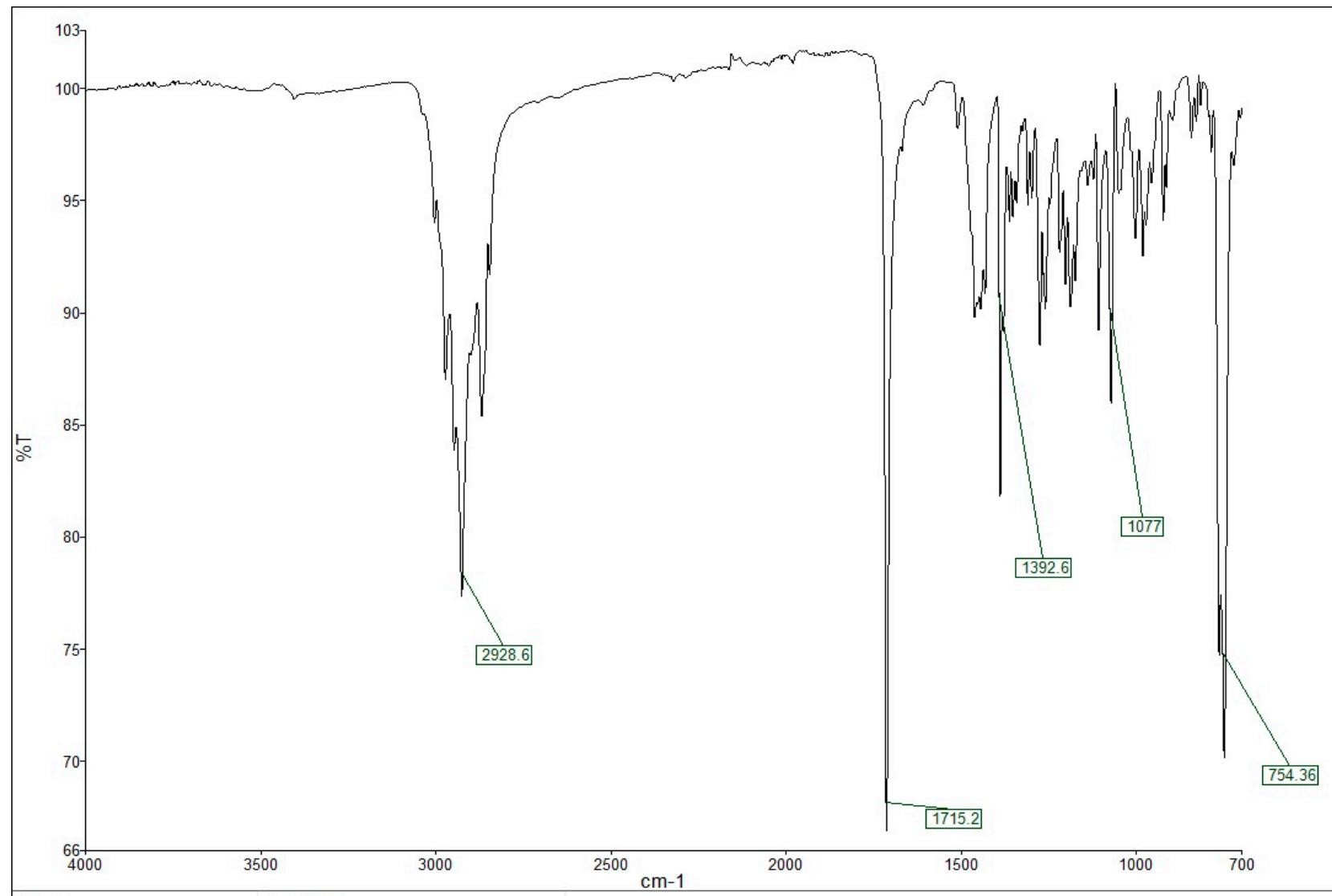


Figure S4: FT-IR spectrum of friedelan-3-one (**1**).

PROTON_05
ZTPA324-495_2_-

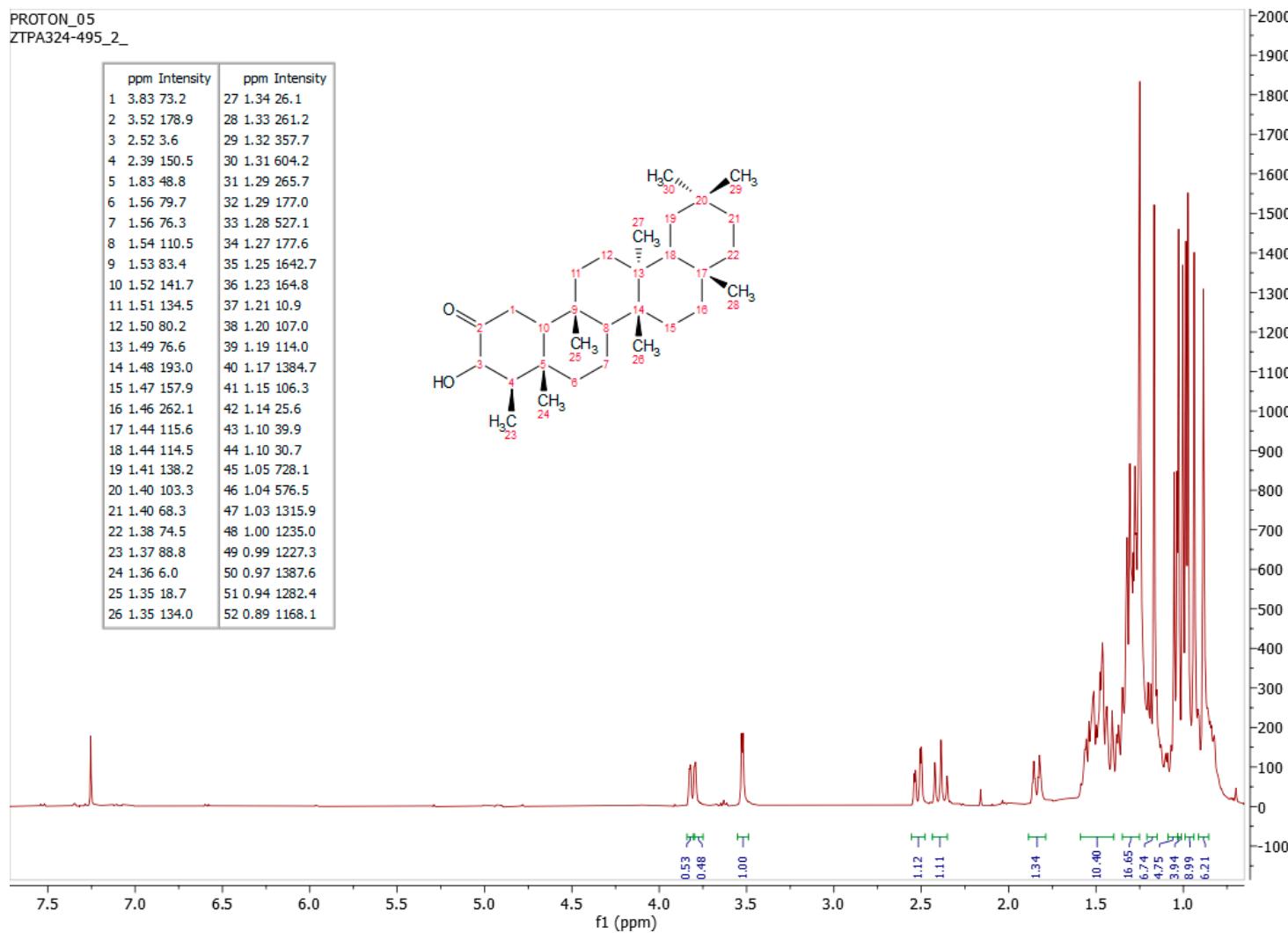


Figure S5: ^1H NMR spectrum of 3α -Hydroxyfriedel-2-one (**2**) recorded in CDCl_3 .

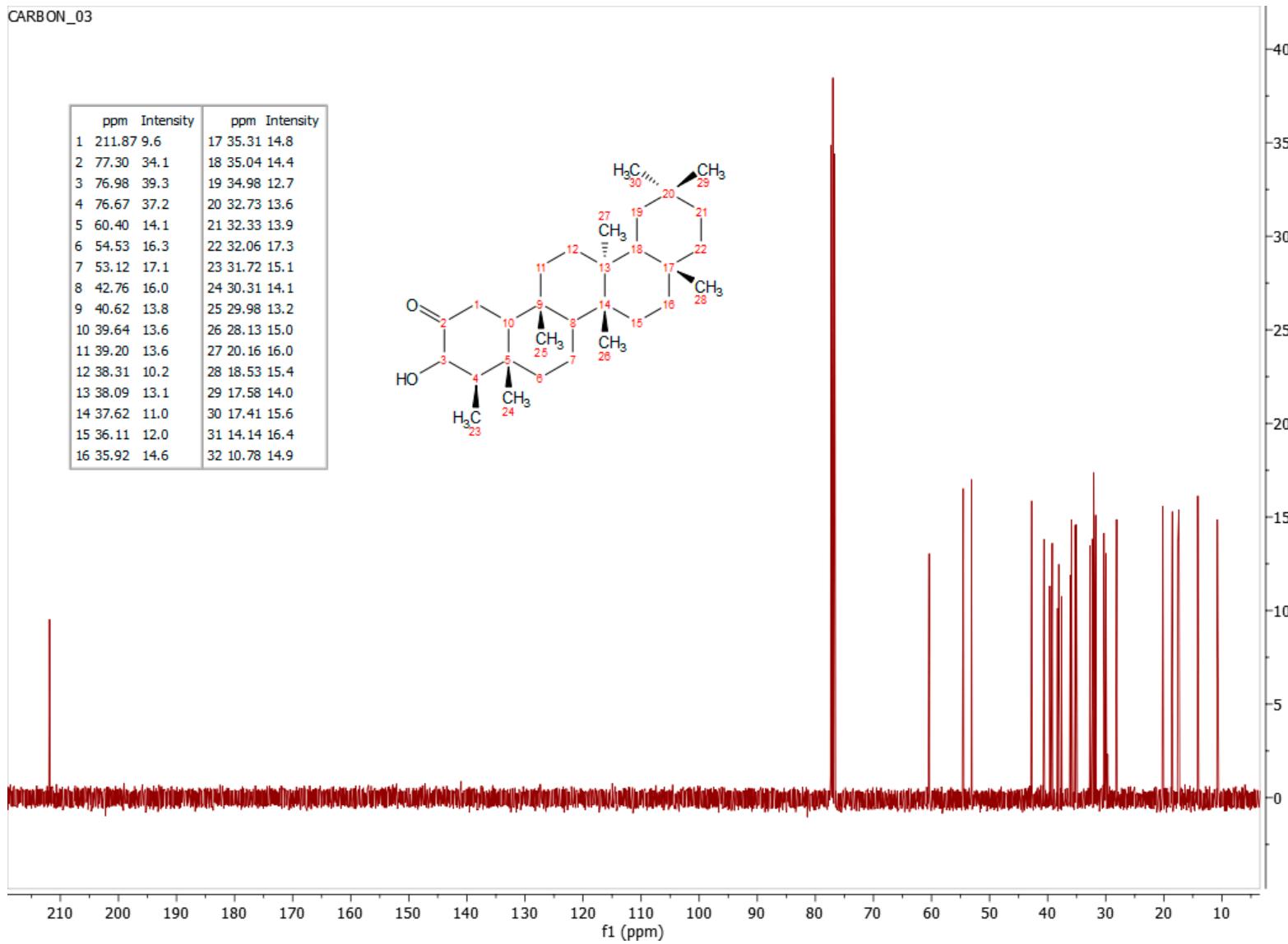


Figure S6: ¹³C NMR spectrum of 3 α -Hydroxyfriedel-2-one (**2**) recorded in CDCl₃.

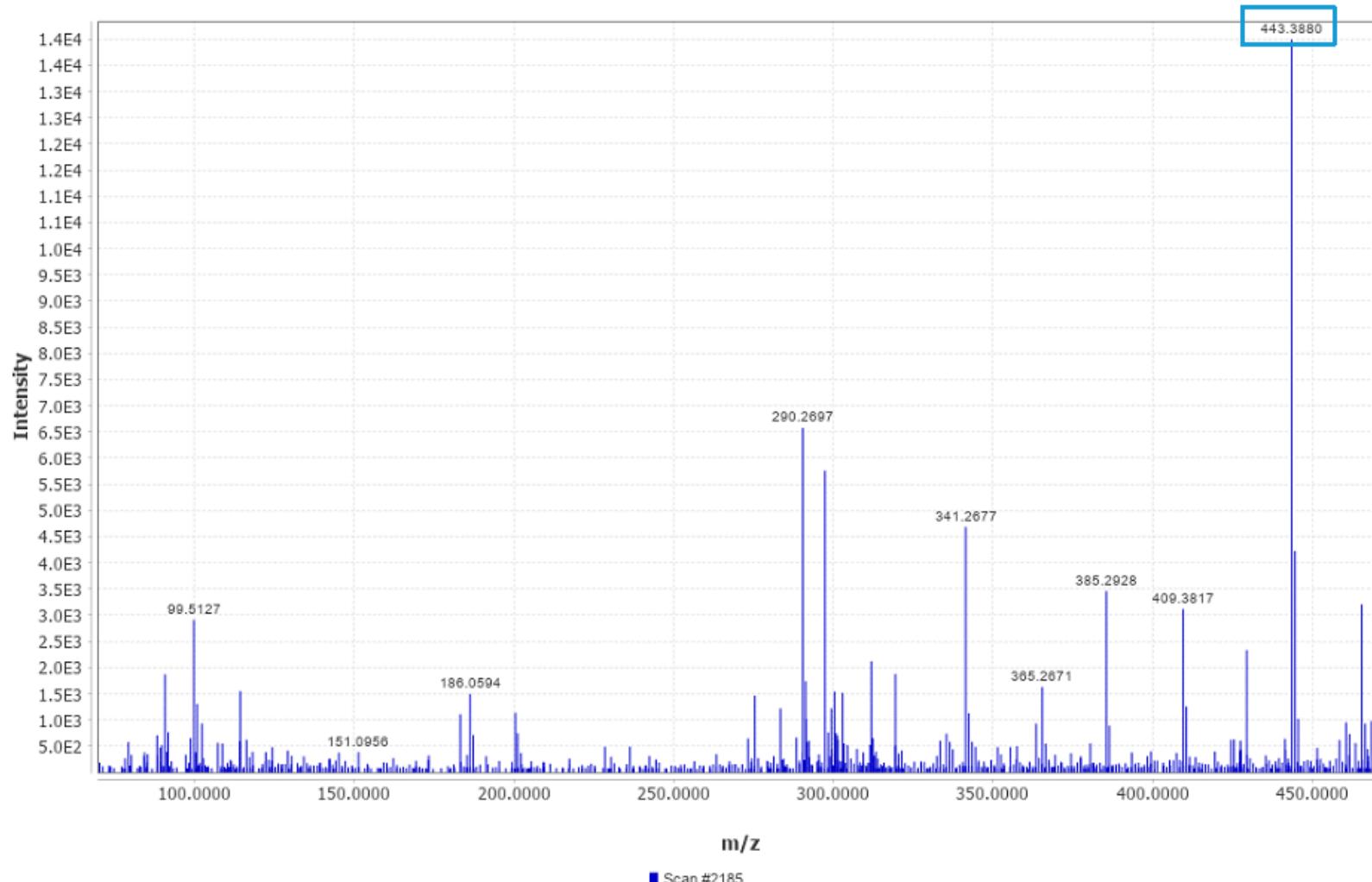


Figure S7: MS spectrum of 3α -Hydroxyfriedel-2-one (**2**).

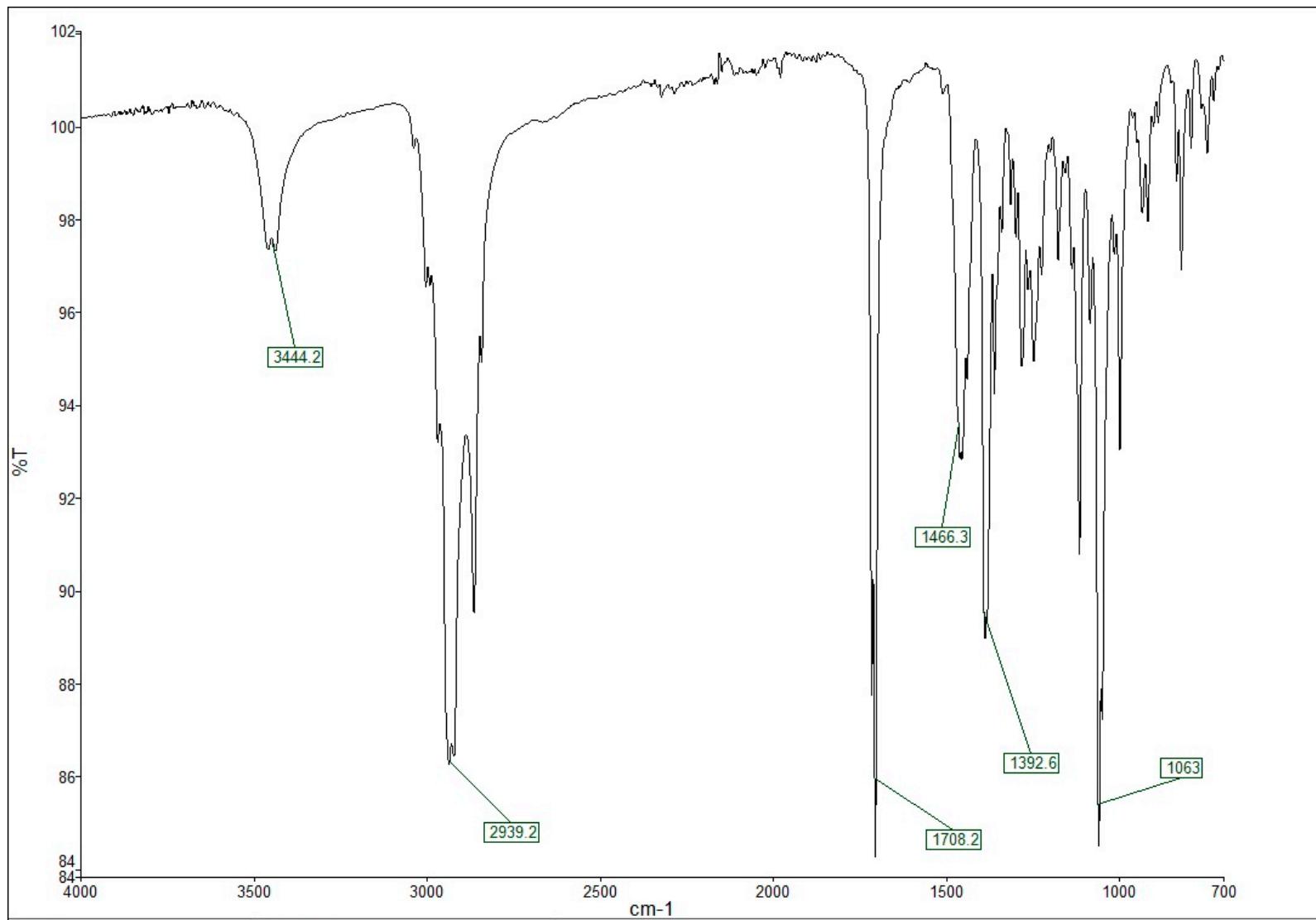


Figure S8: FT-IR spectrum of 3 α -Hydroxyfriedel-2-one (**2**).

PROTON_05
ZTPA318-329_1_cult

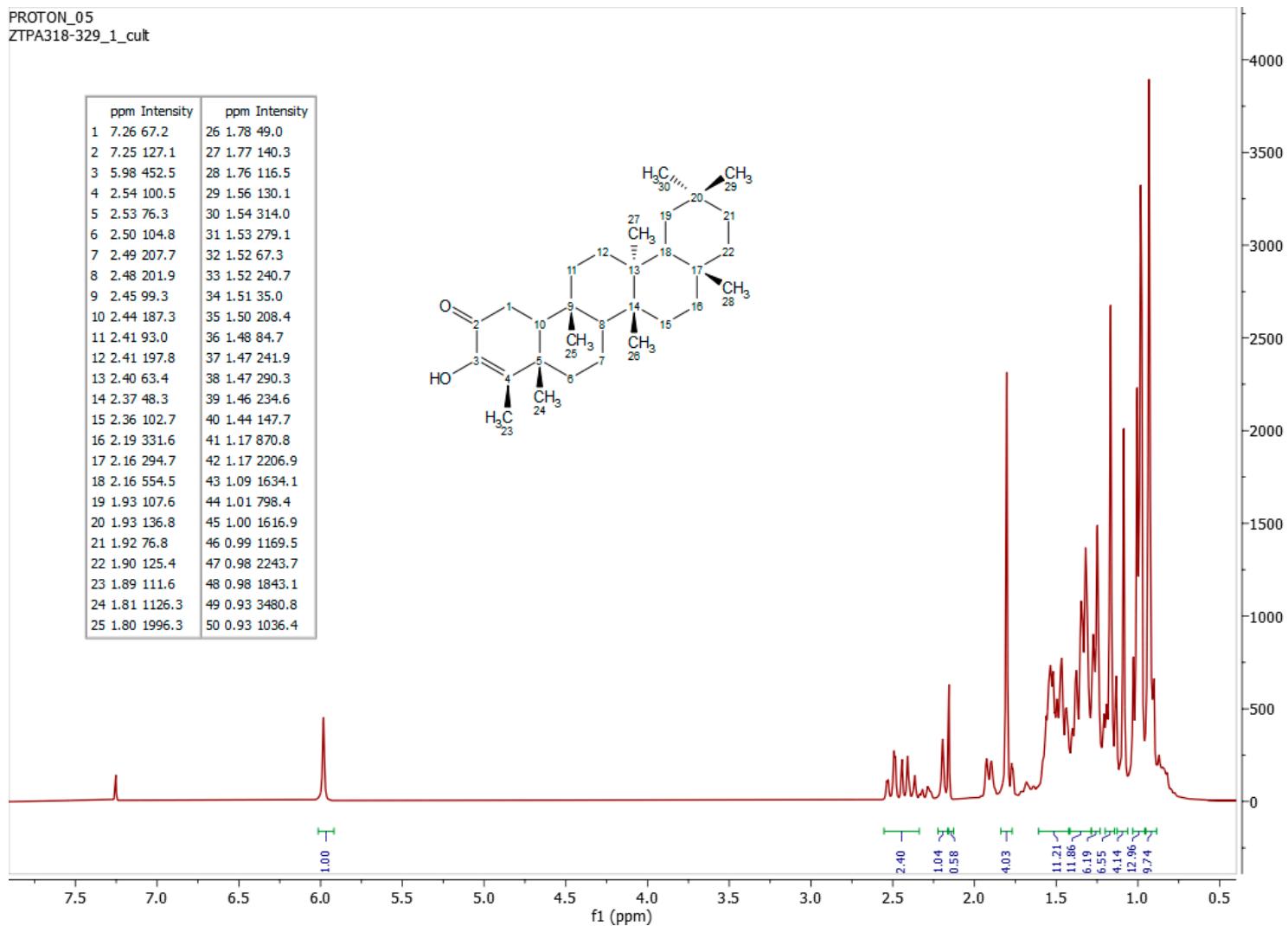


Figure S9: ¹H NMR spectrum of 3-Hydroxyfriedel-3-en-2-one (**3**) recorded in CDCl₃.

CARBON_05
ZTPA318-329_1_cult

	ppm	Intensity
1	195.04	119.7
2	142.57	121.2
3	140.79	111.7
4	77.31	430.3
5	76.99	457.4
6	76.68	464.6
7	55.68	200.3
8	52.63	198.0
9	42.77	198.3
10	39.71	171.8
11	39.53	154.8
12	39.20	193.3
13	38.39	193.7
14	38.24	215.6
15	36.75	188.8
16	35.90	160.1
	32	10.37 183.5
	17	35.30 155.8
	18	34.96 181.8
	19	34.69 169.8
	20	32.74 146.1
	21	32.14 161.5
	22	32.08 167.9
	23	31.77 193.5
	24	30.22 147.5
	25	29.99 180.9
	26	28.13 207.6
	27	20.07 199.6
	28	18.87 169.8
	29	18.50 248.6
	30	17.97 143.9
	31	17.66 208.8
	32	10.37 183.5

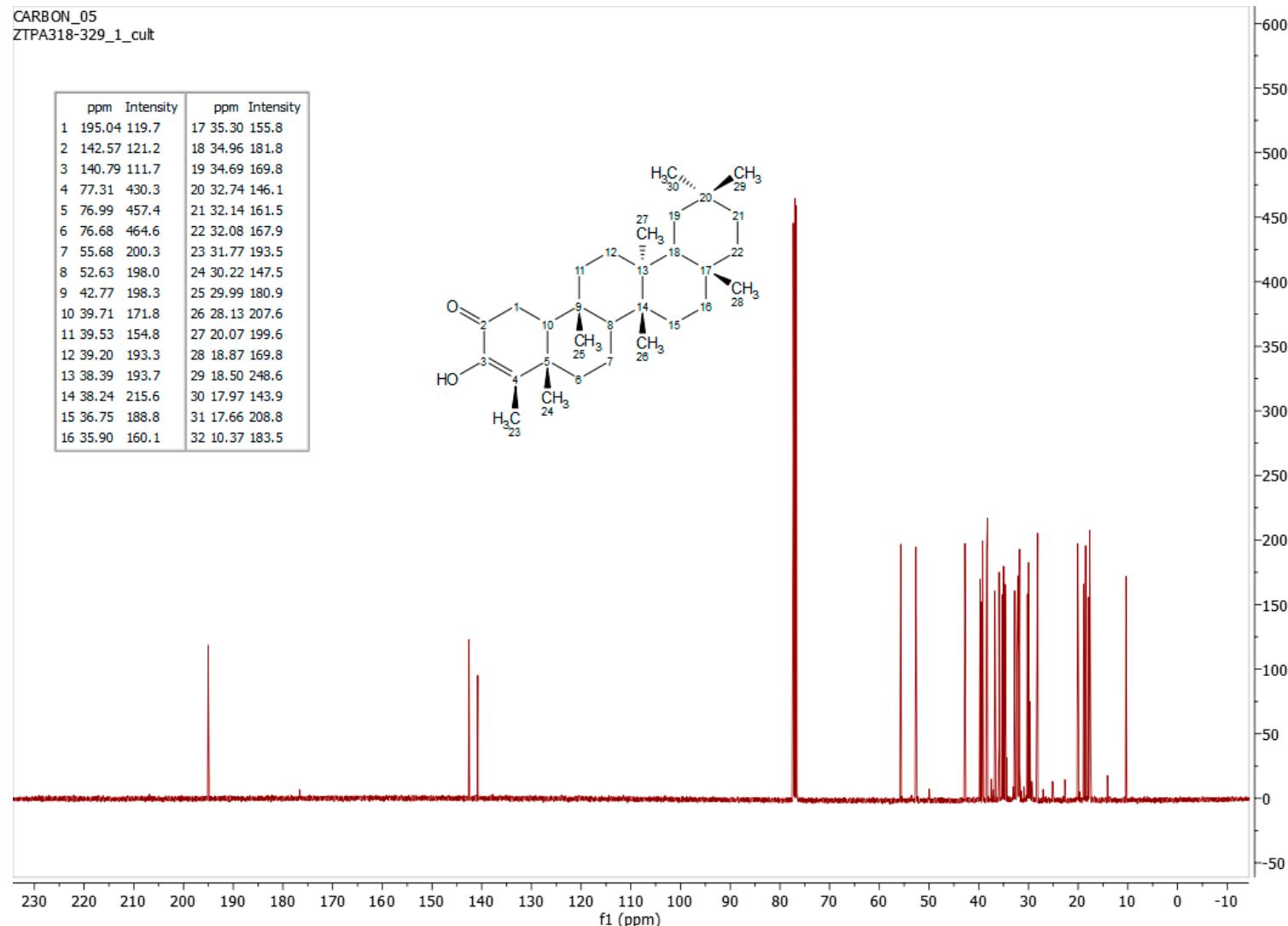
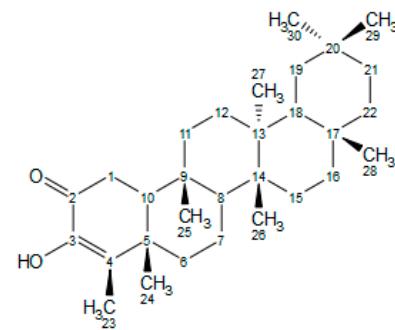


Figure S10: ¹³C NMR spectrum of 3-Hydroxyfriedel-3-en-2-one (**3**) recorded in CDCl₃.

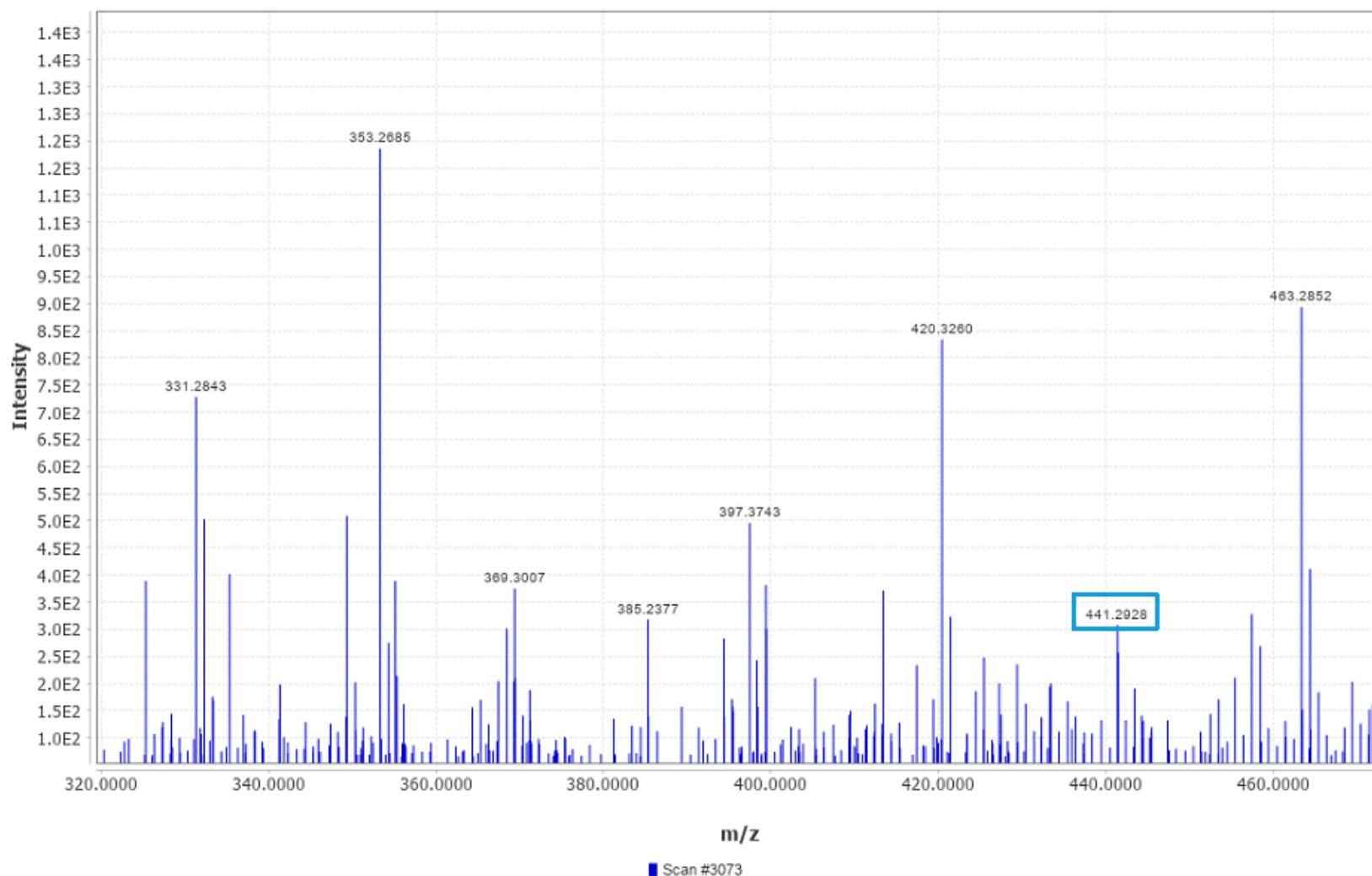


Figure S11: MS spectrum of 3-Hydroxyfriedel-3-en-2-one (**3**).

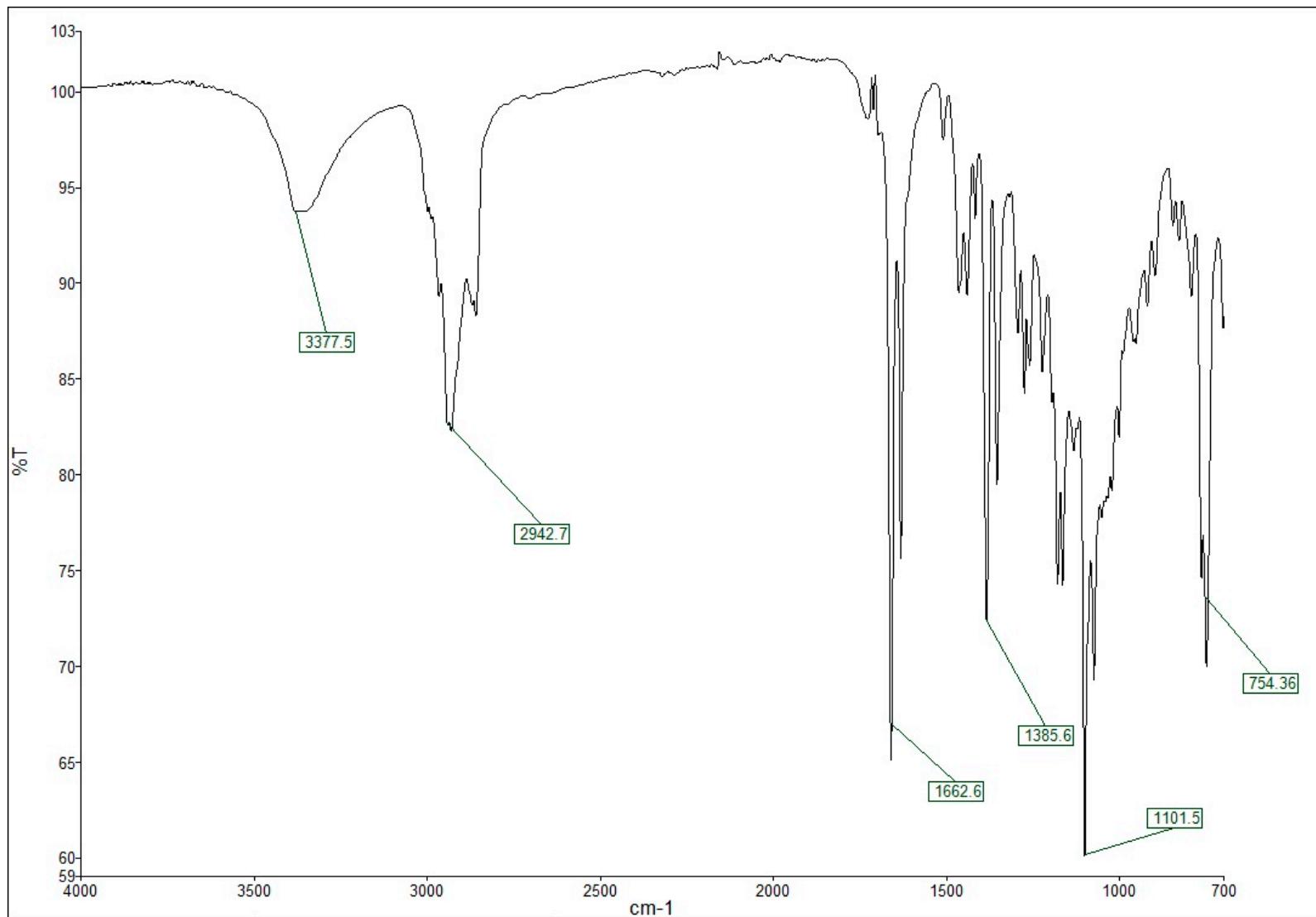


Figure S12: FT-IR spectrum of 3-Hydroxyfriedel-3-en-2-one (**3**)

PROTON_05
ZTPA203-237_X2_cult

	ppm Intensity	ppm Intensity
1	4.68 293.3	26 1.37 93.7
2	4.56 296.0	27 1.36 491.0
3	3.19 142.9	28 1.35 81.9
4	2.38 114.3	29 1.33 60.4
5	1.91 75.7	30 1.33 183.3
6	1.67 1604.4	31 1.29 90.4
7	1.63 117.3	32 1.28 73.0
8	1.61 111.2	33 1.26 452.2
9	1.59 181.0	34 1.25 870.1
10	1.58 134.7	35 1.22 151.3
11	1.52 113.9	36 1.20 133.8
12	1.50 105.3	37 1.20 129.6
13	1.49 108.4	38 1.17 128.4
14	1.47 91.8	39 1.13 147.0
15	1.46 114.9	40 1.13 152.1
16	1.43 77.1	41 1.02 2244.6
17	1.42 120.3	42 0.96 2363.8
18	1.42 110.3	43 0.94 1106.7
19	1.41 176.7	44 0.94 1410.2
20	1.40 119.9	45 0.82 1462.8
21	1.40 131.7	46 0.82 887.2
22	1.39 237.8	47 0.78 1935.0
23	1.38 658.2	48 0.75 2291.3
24	1.38 443.2	49 0.69 76.2
25	1.37 74.4	50 0.69 90.2

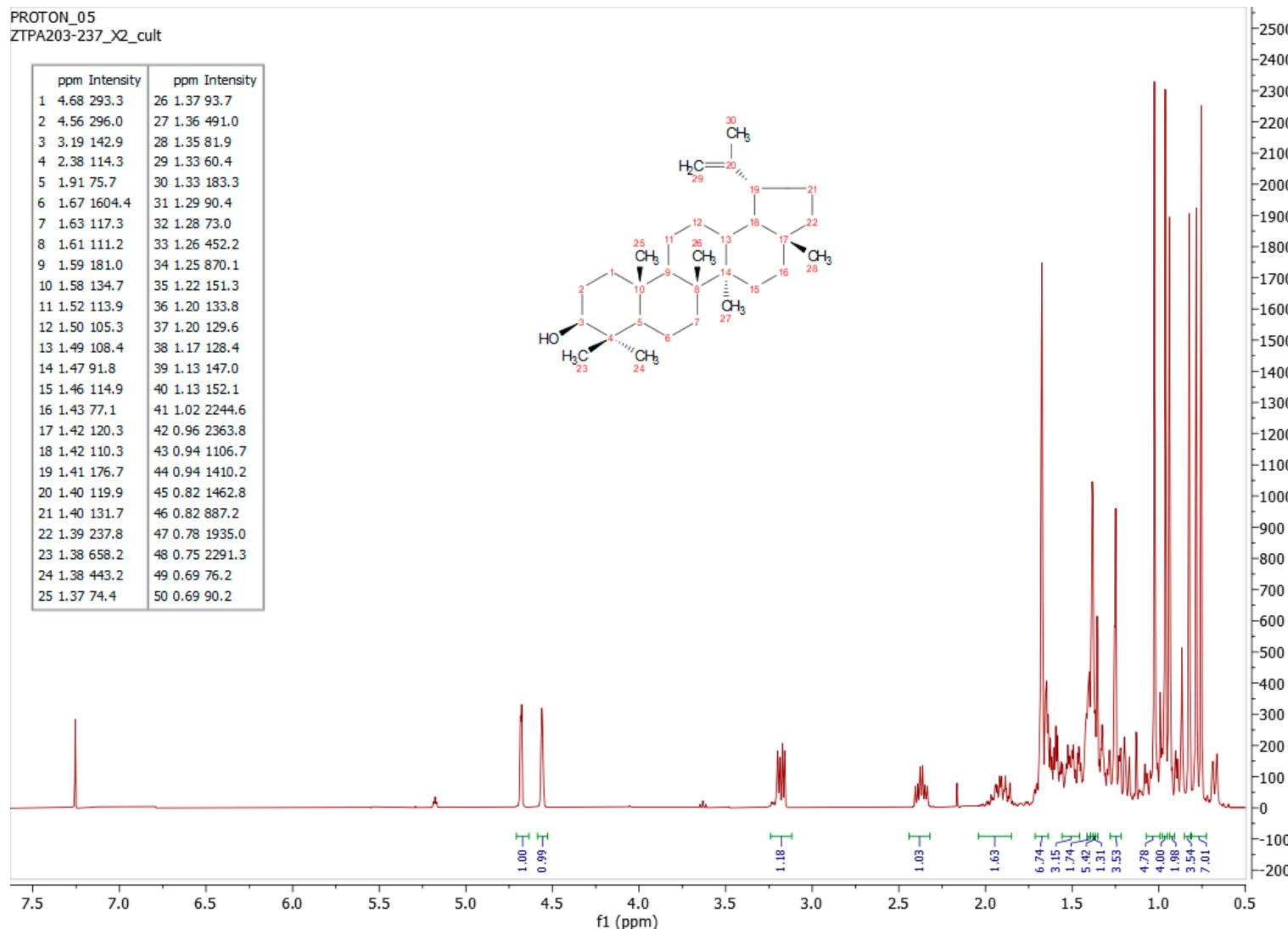
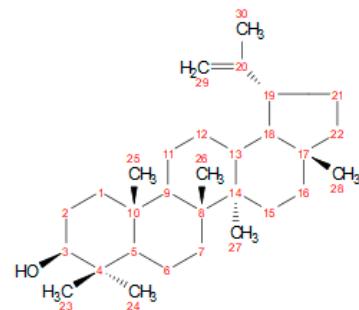


Figure S13: ^1H NMR spectrum of lupeol (**4**) recorded in CDCl_3 .

CARBON_04
ZTPA203-237_X2_cult

ppm	Intensity	ppm	Intensity		
1	150.96	104.8	18	38.03	160.7
2	109.31	43.4	19	37.15	139.4
3	79.02	16.2	20	35.57	101.5
4	78.99	161.2	21	34.26	101.3
5	77.32	481.7	22	29.83	107.6
6	77.01	475.7	23	29.69	53.5
7	76.69	482.3	24	27.98	127.7
8	55.28	161.8	25	27.43	111.1
9	50.42	144.0	26	27.39	108.0
10	48.28	134.9	27	25.12	103.5
11	47.97	95.3	28	20.91	102.6
12	42.99	156.0	29	18.31	98.6
13	42.82	123.1	30	17.99	161.4
14	40.81	132.8	31	16.11	178.8
15	39.99	99.7	32	15.96	89.1
16	38.85	151.4	33	15.36	157.3
17	38.69	131.0	34	14.54	114.1

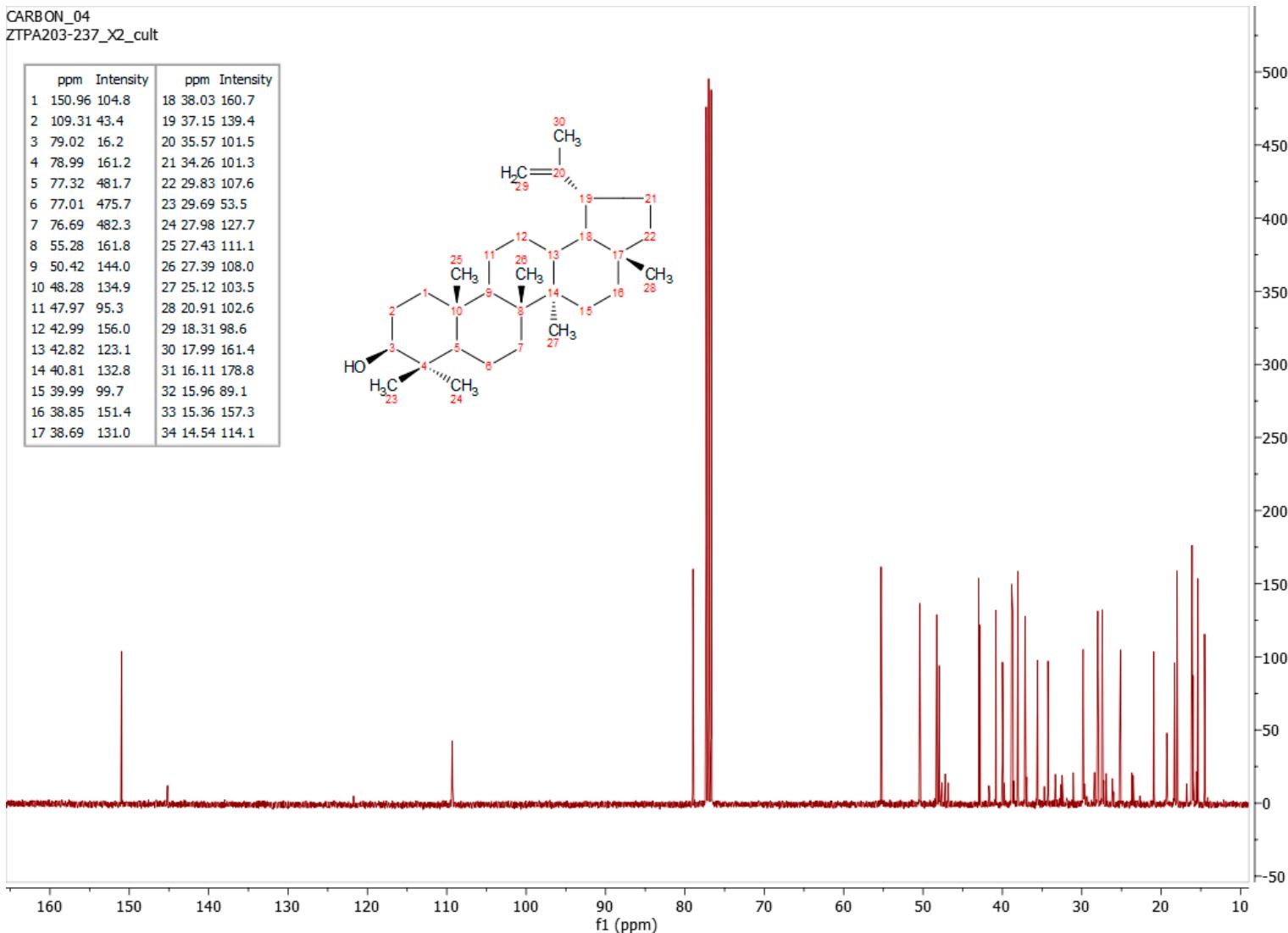
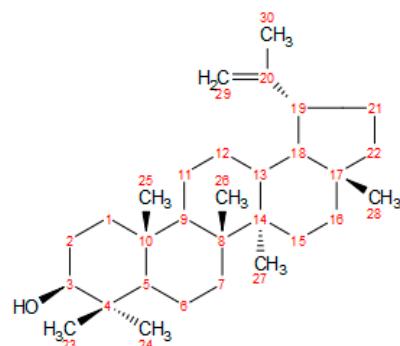


Figure S14: ^{13}C NMR spectrum of lupeol (**4**) recorded in CDCl_3 .

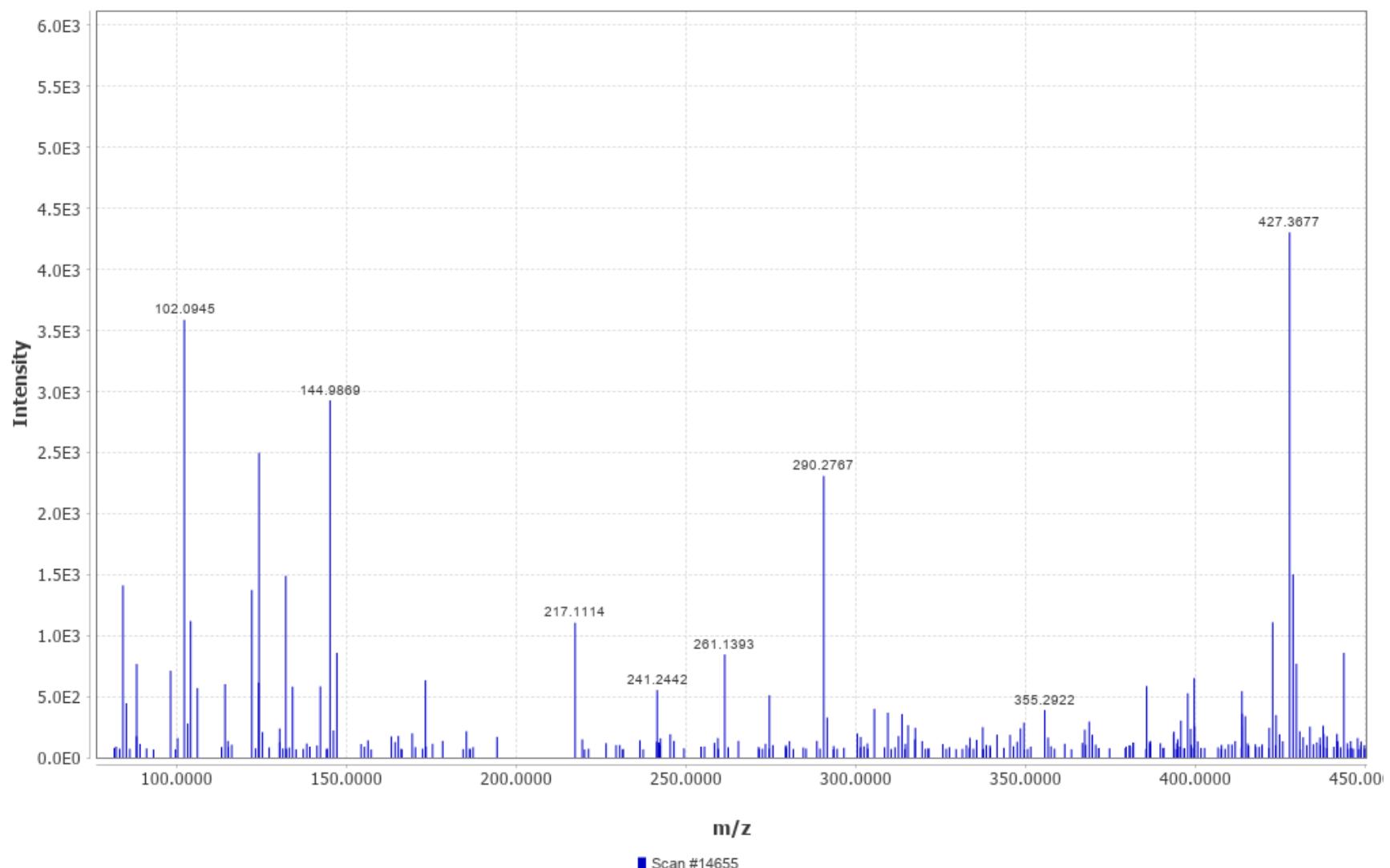


Figure S15: MS spectrum of lupeol (4).

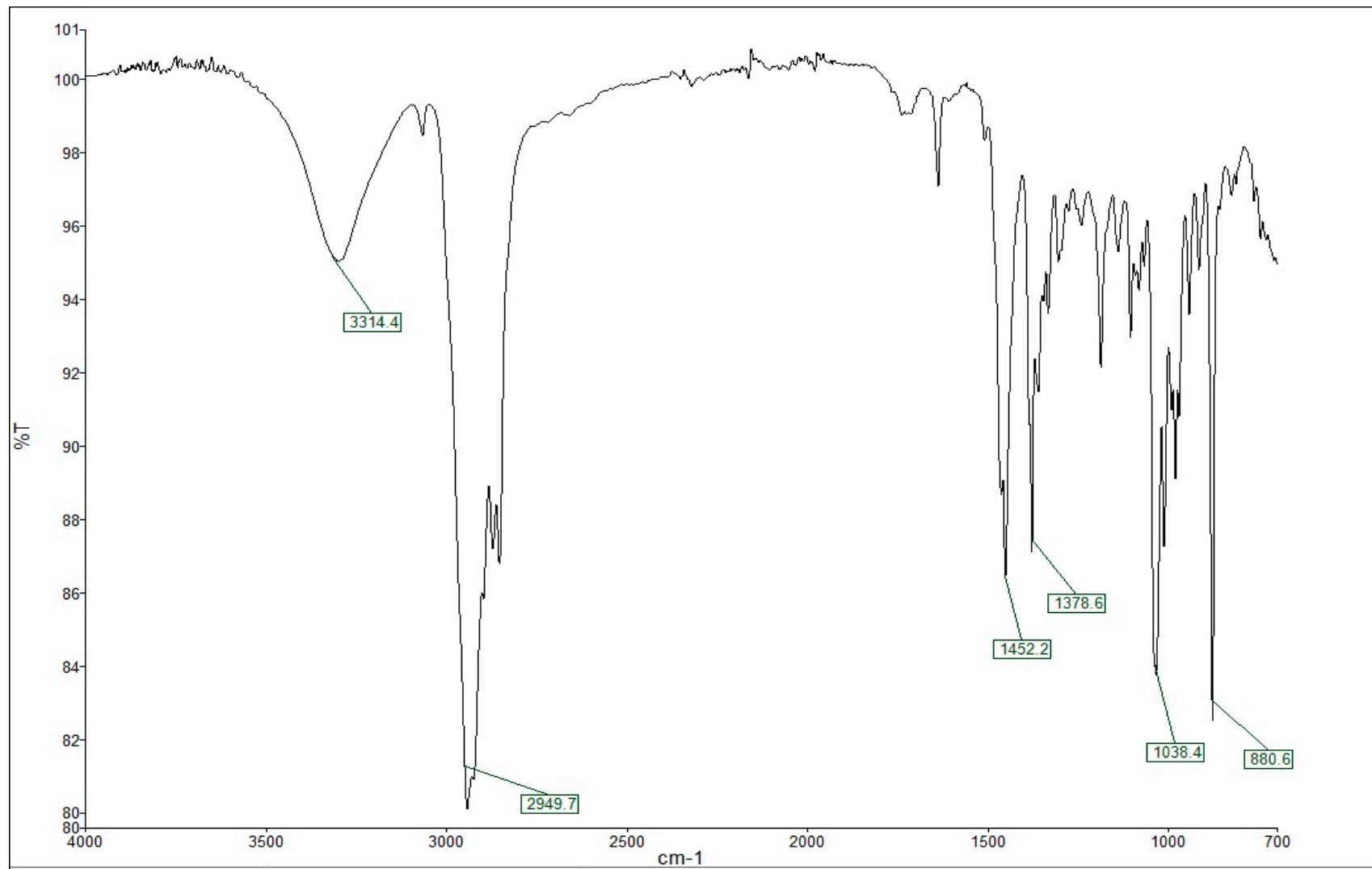


Figure S16: FT-IR spectrum of lupeol (**4**)

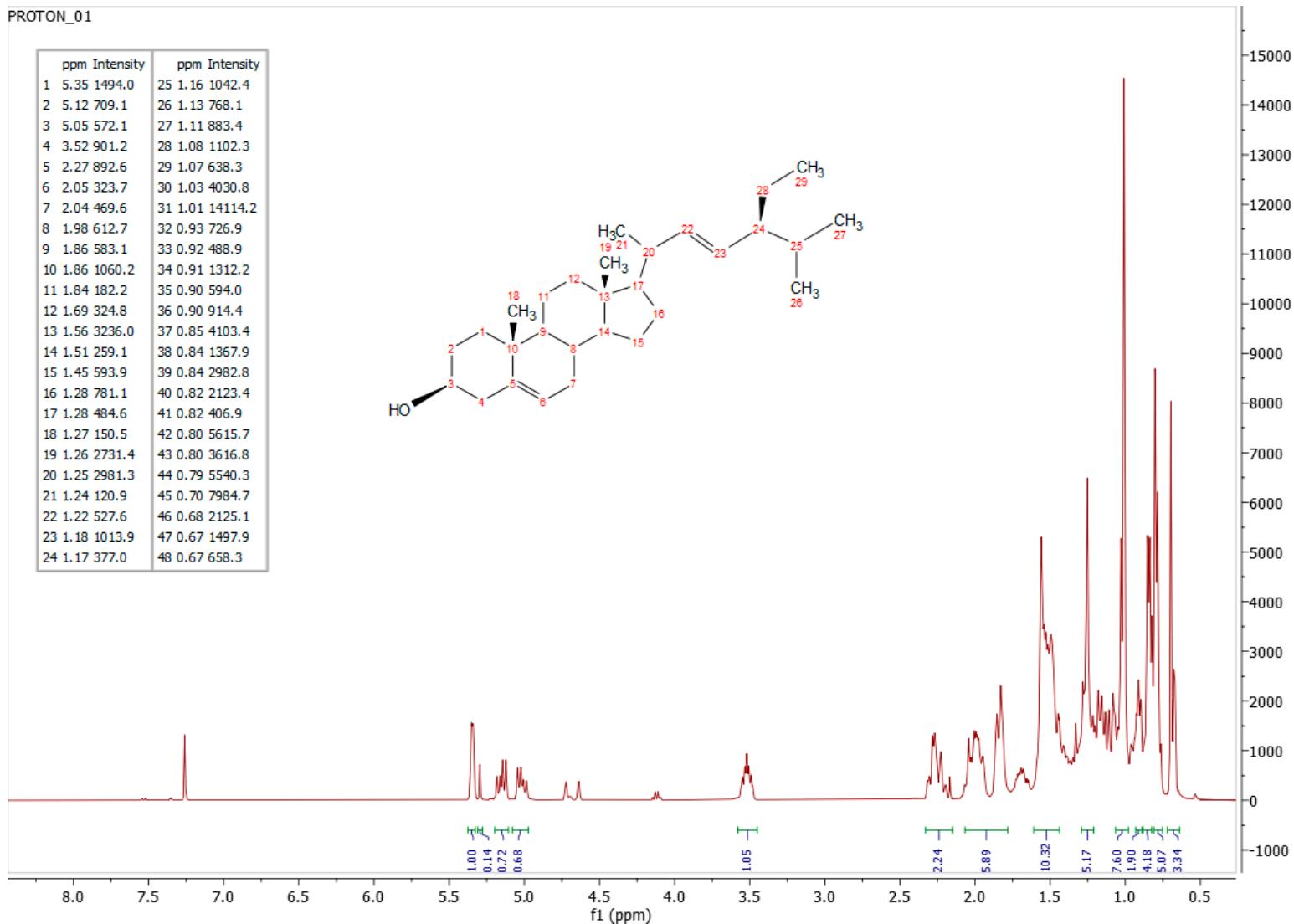


Figure S17: ^1H NMR spectrum of stigmasterol (**5**) recorded in CDCl_3 .

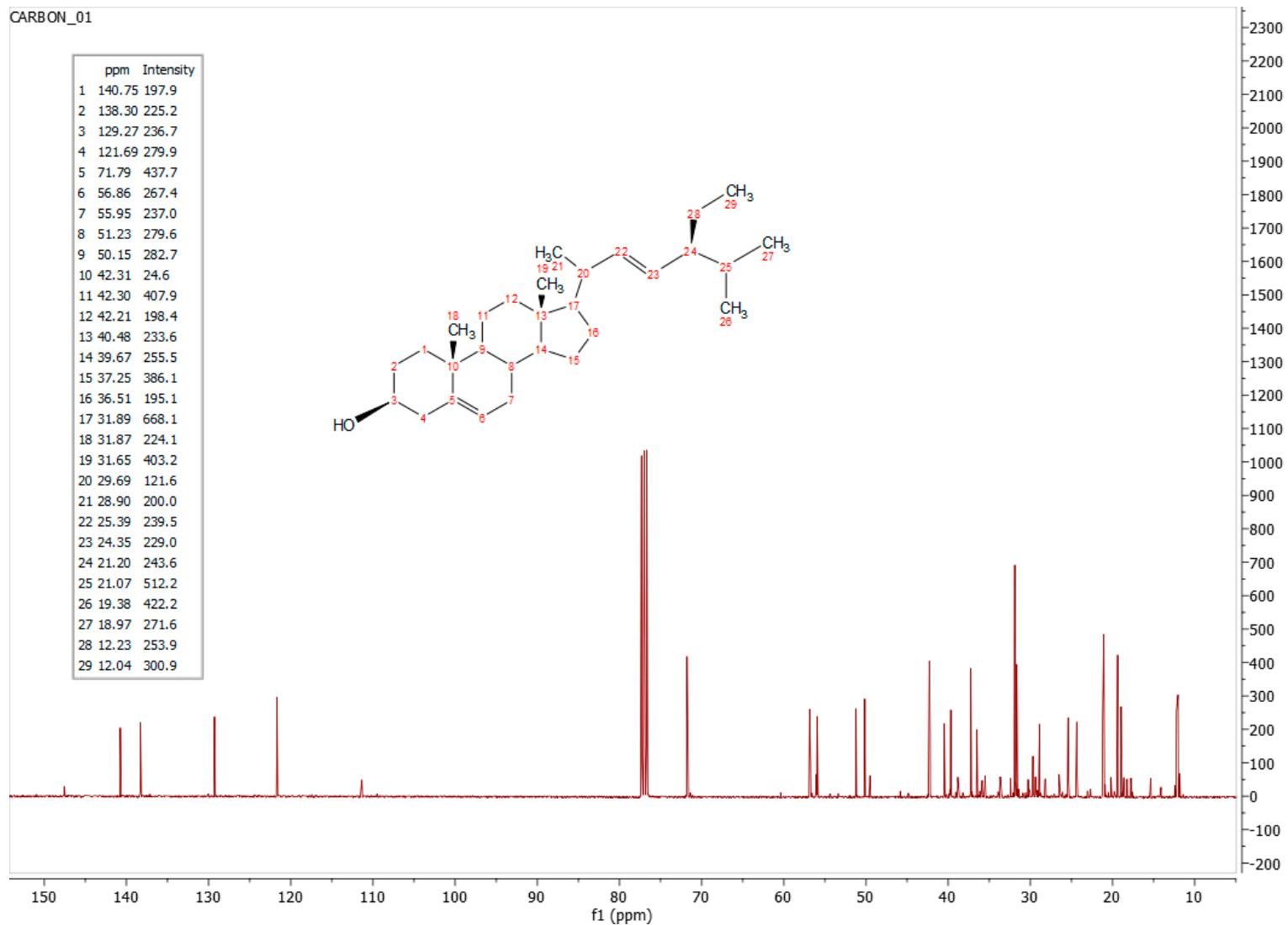


Figure S18: ^{13}C NMR spectrum of stigmasterol (**5**) recorded in CDCl_3

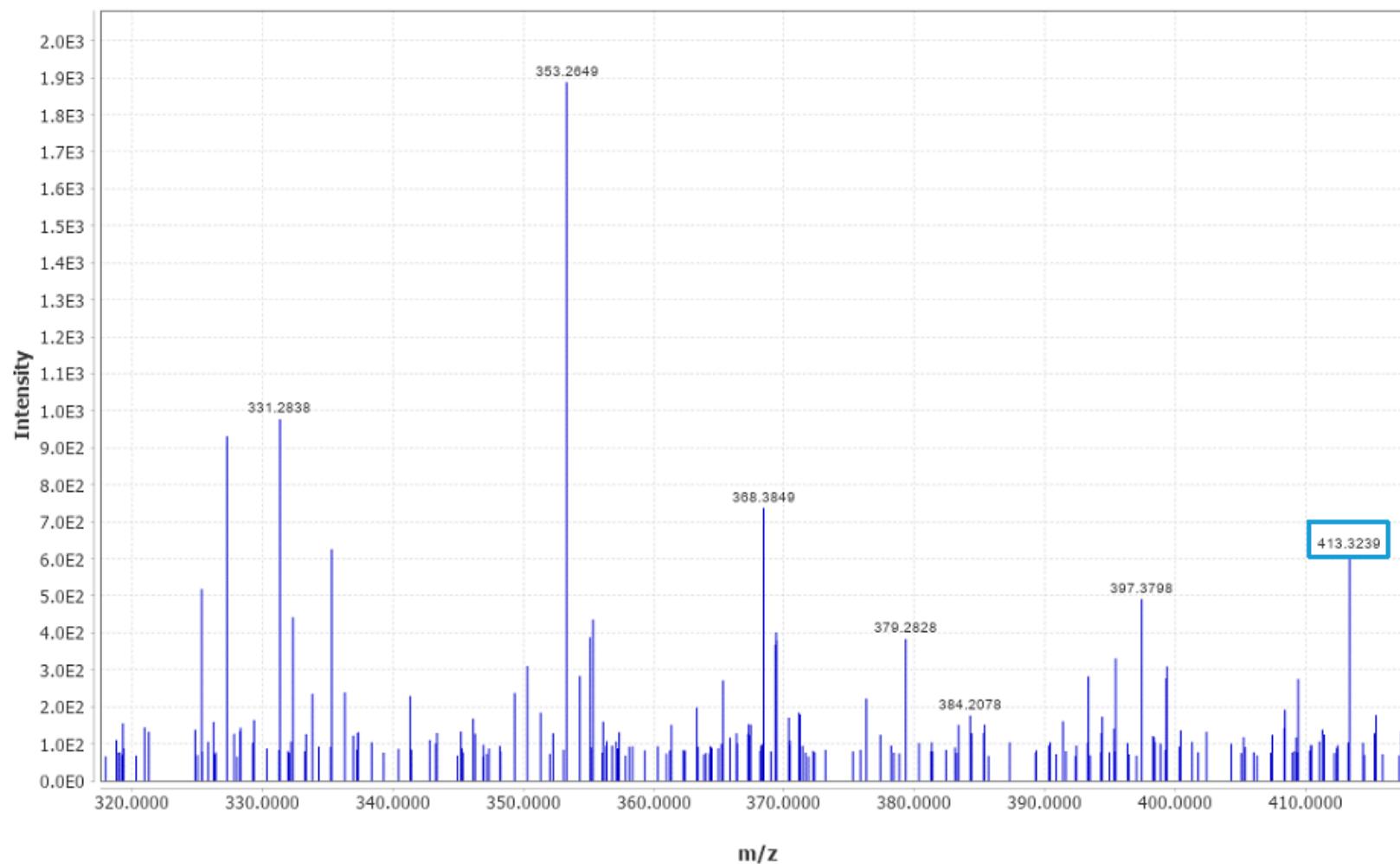


Figure S19: MS spectrum of stigmasterol (**5**).

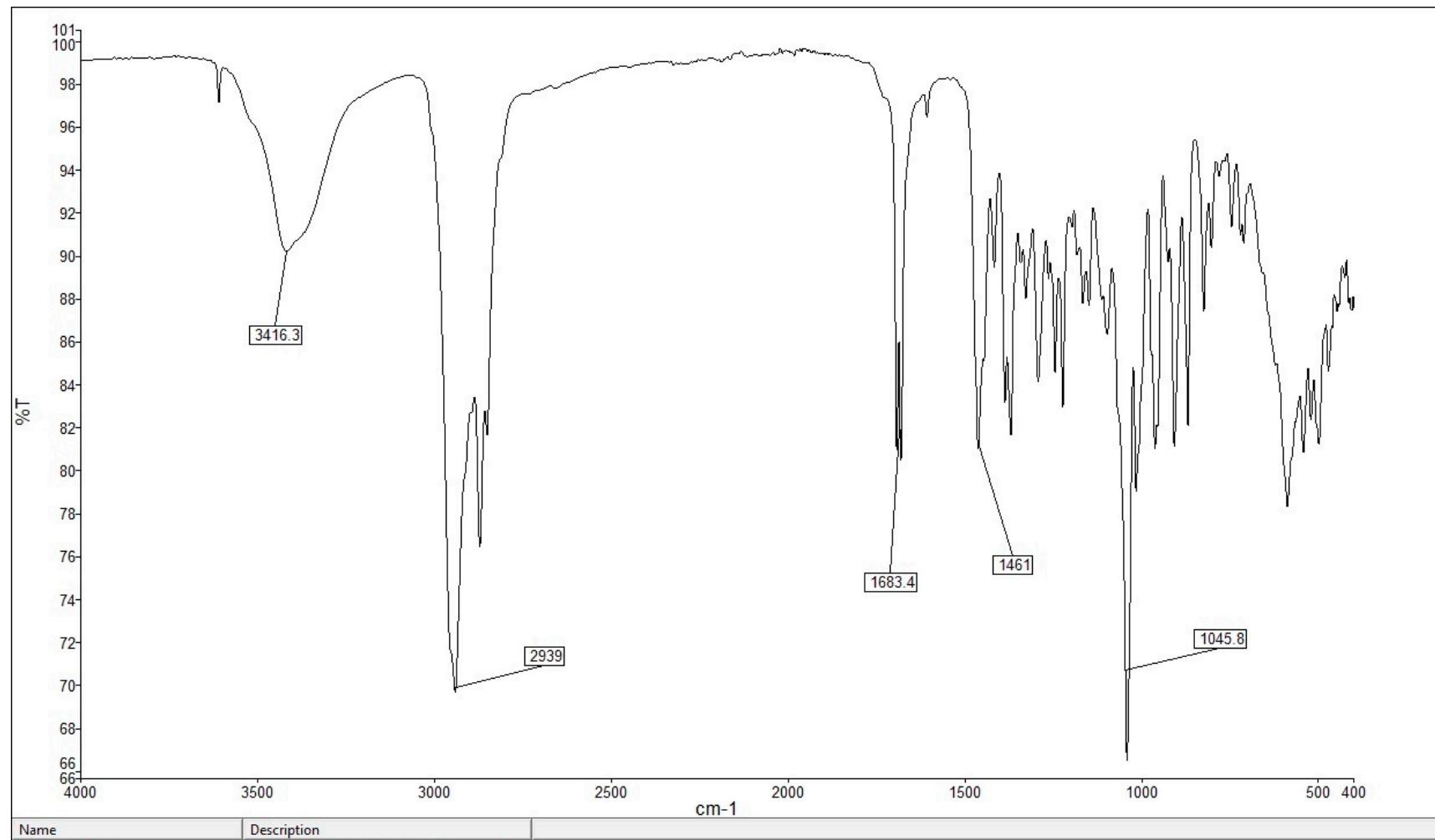


Figure S20: FT-IR spectrum of stigmasterol (**5**)

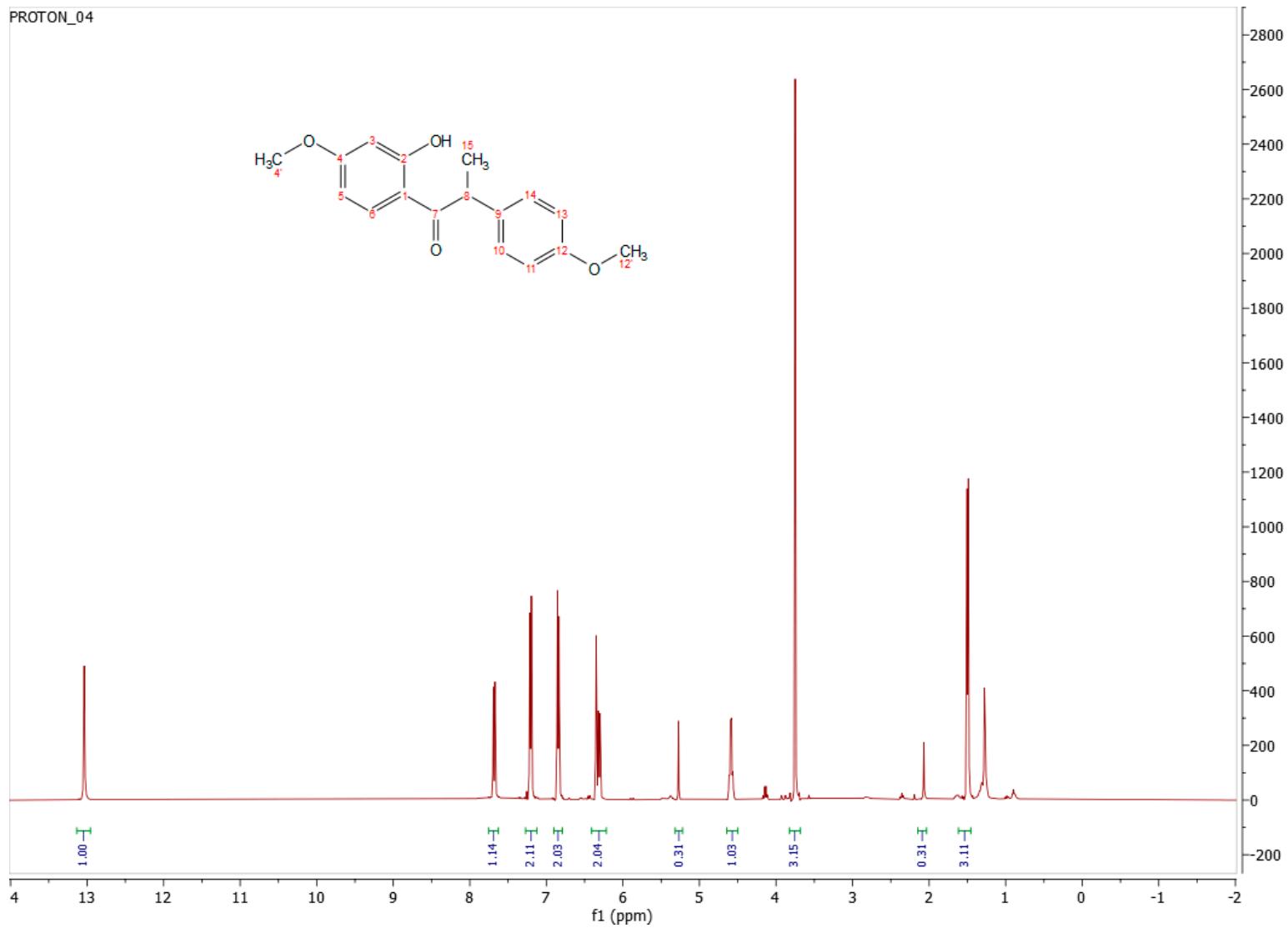


Figure S21: ¹H NMR spectrum of (±)-4-O-Methylangolensin (**6**) recorded in CDCl₃.

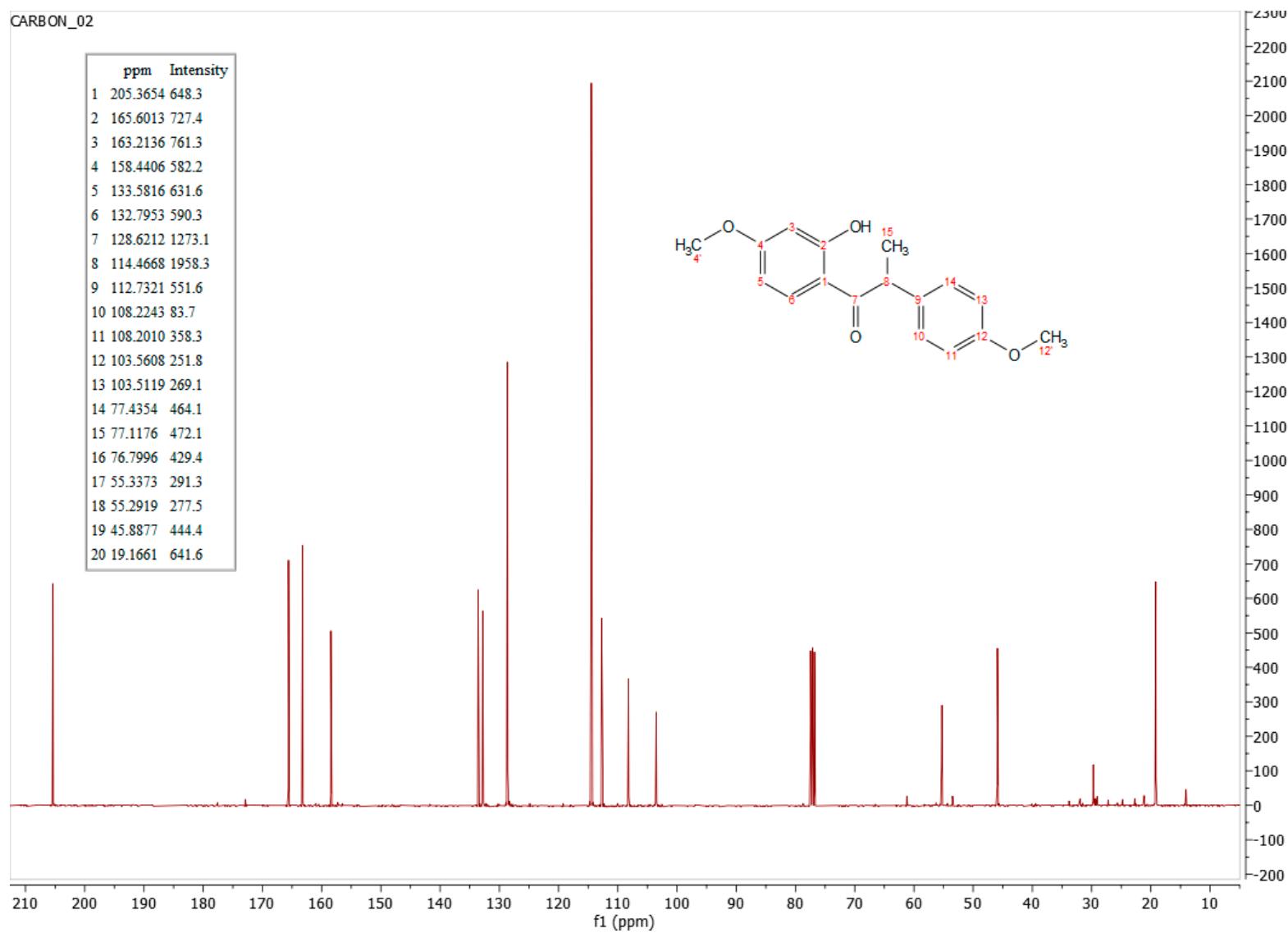


Figure S22: ^{13}C NMR spectrum of (\pm) -4-O-Methylangolensin (**6**) recorded in CDCl_3

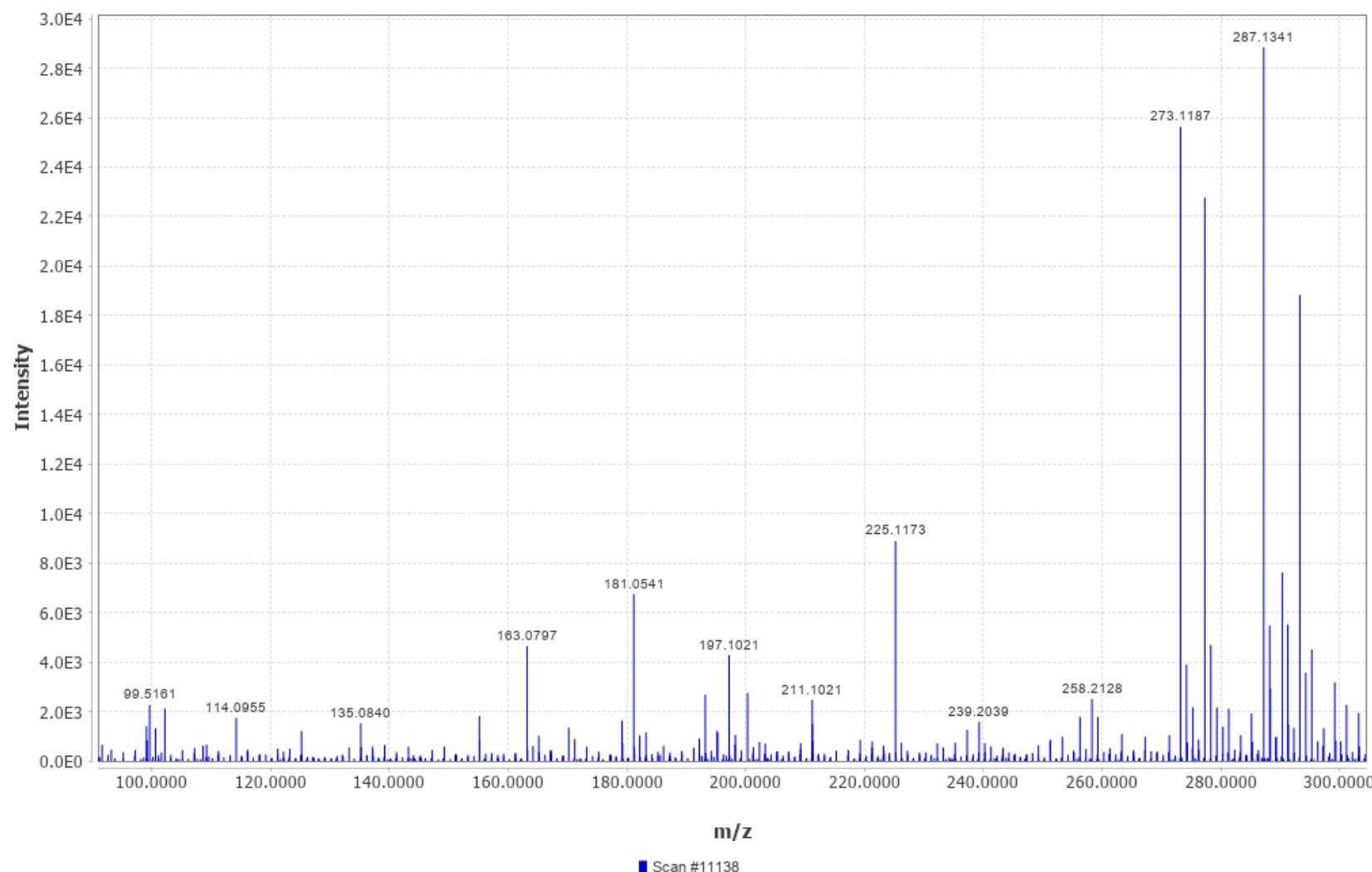


Figure S23: MS spectrum of (\pm) -4-O-Methylangolensin (**6**).

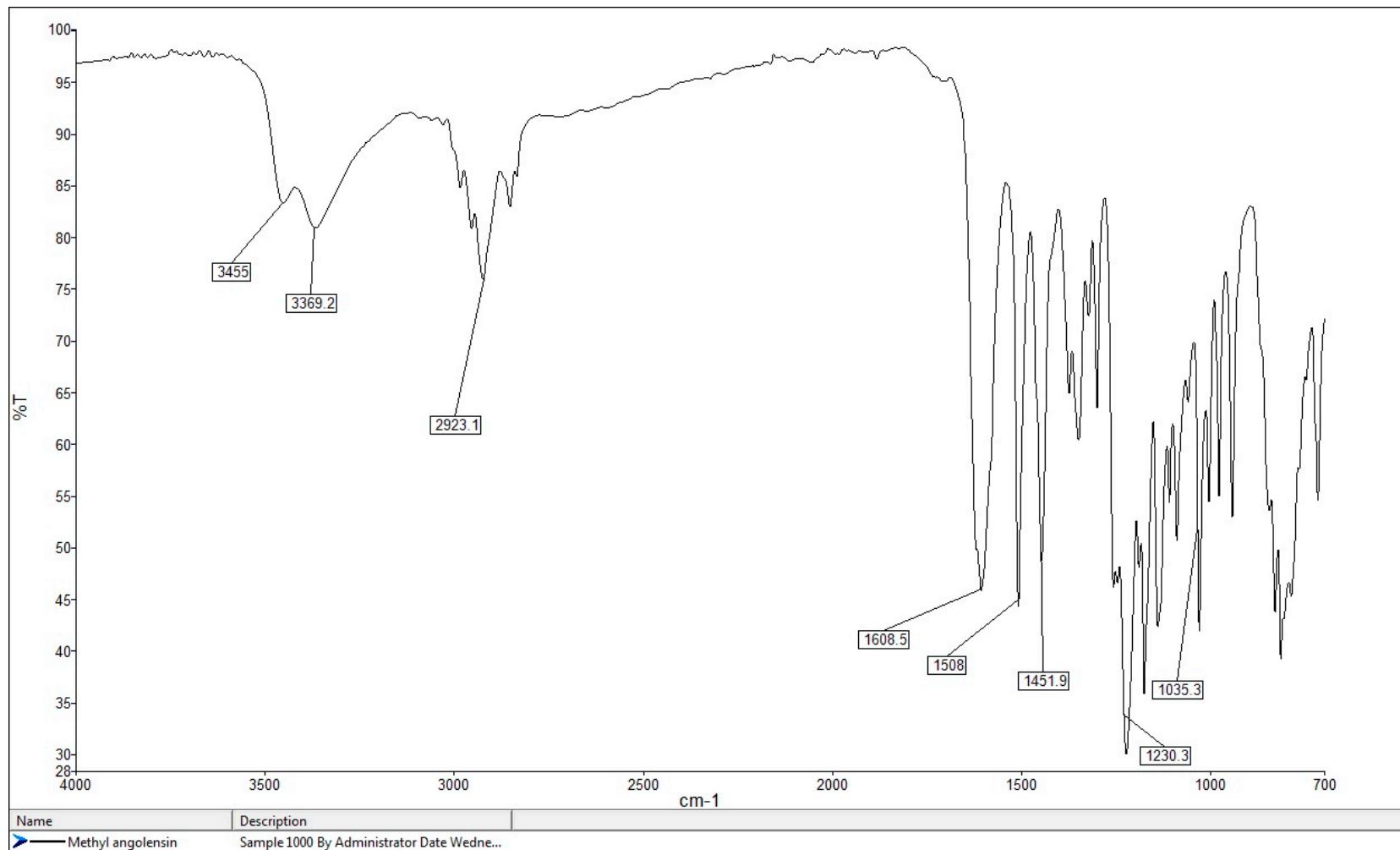


Figure S24: FT-IR spectrum of (\pm) -4-O-Methylangolensin (7).

PROTON_02
ZTPA 86-106_light

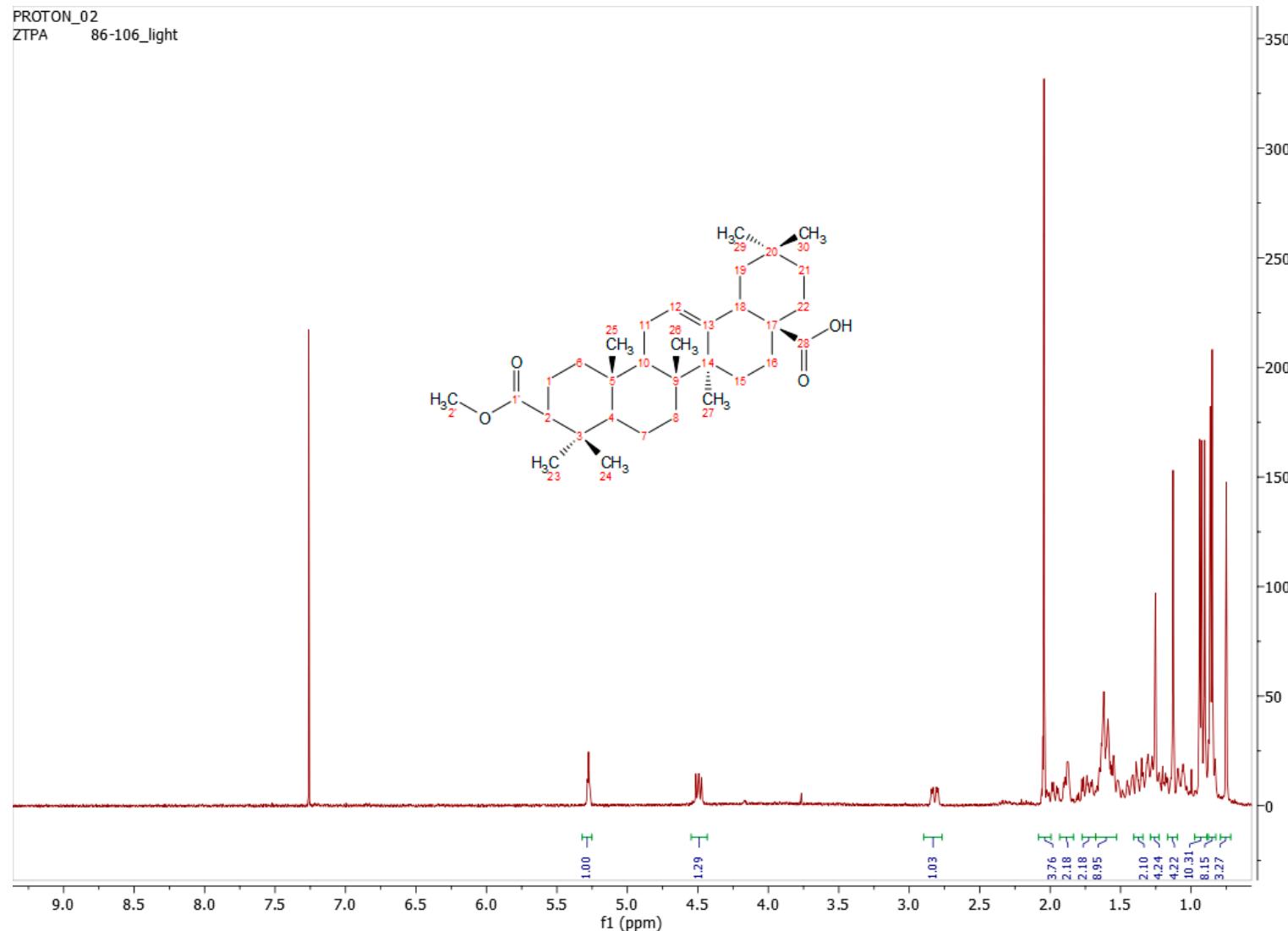


Figure S25: ¹H NMR spectrum of Oleanolic acid acetate (7) recorded in CDCl₃.

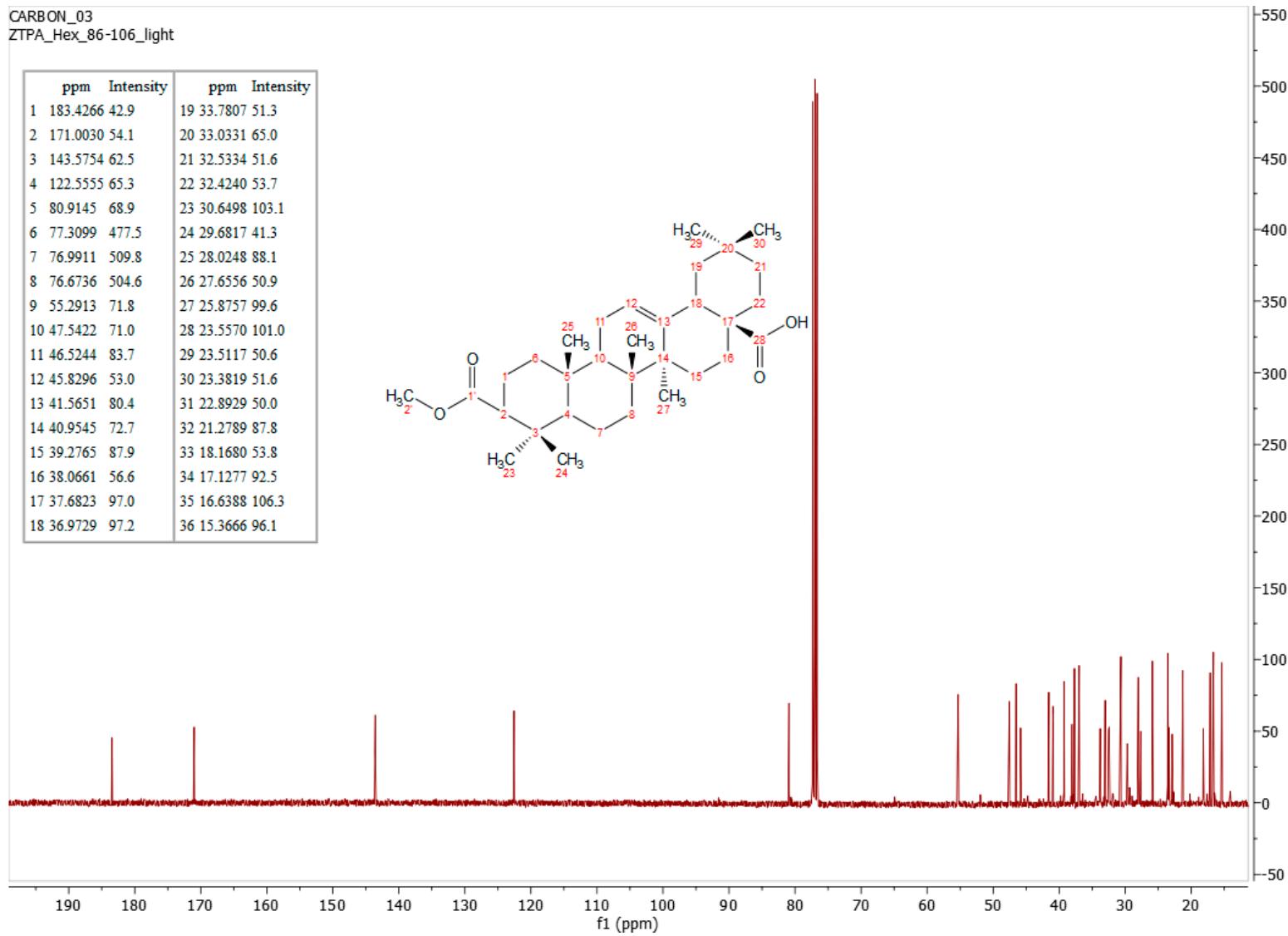


Figure S26: ¹³C NMR spectrum of Oleanolic acid acetate (**7**) recorded in CDCl₃

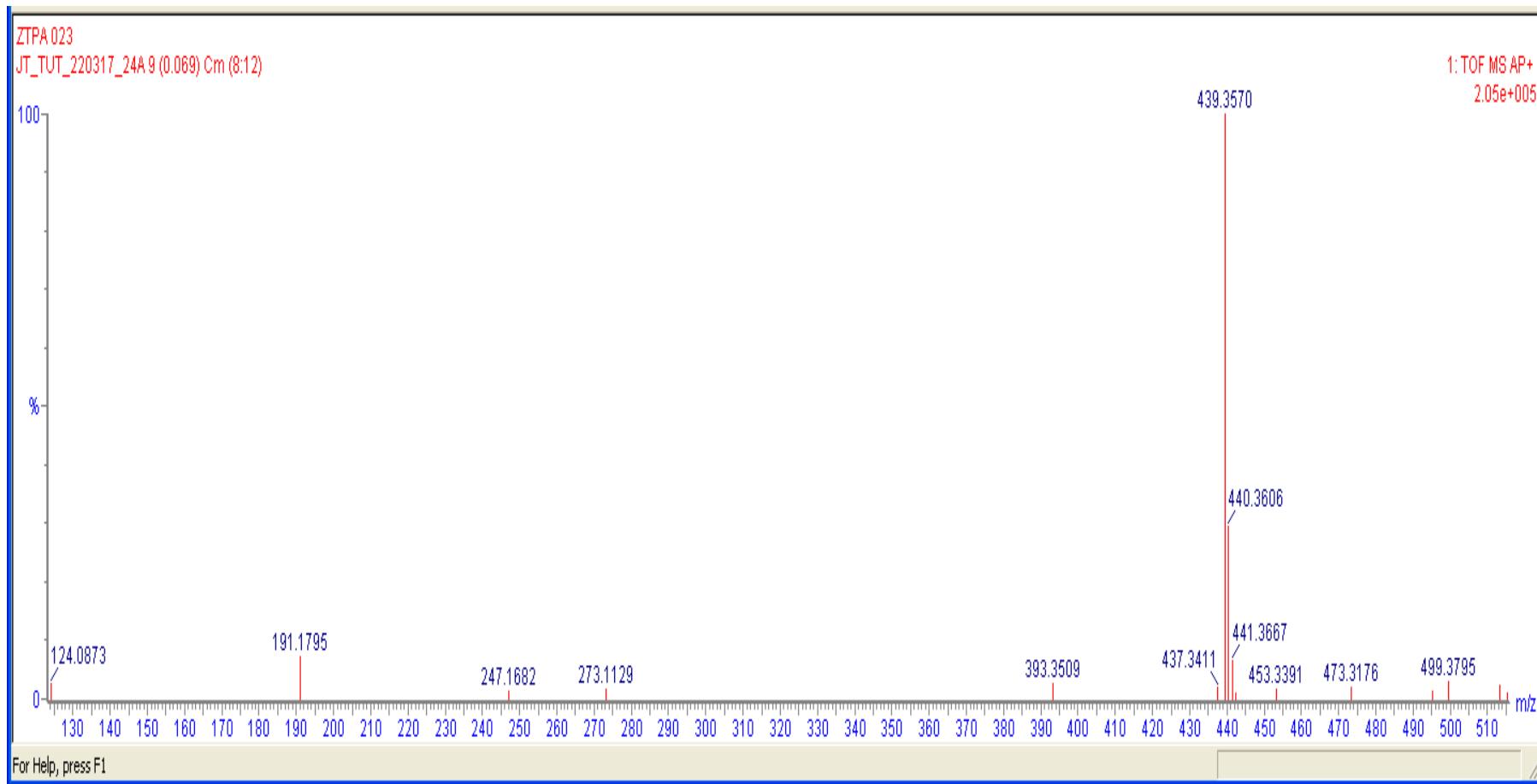


Figure S27: MS spectrum of Oleanolic acid acetate (**7**).

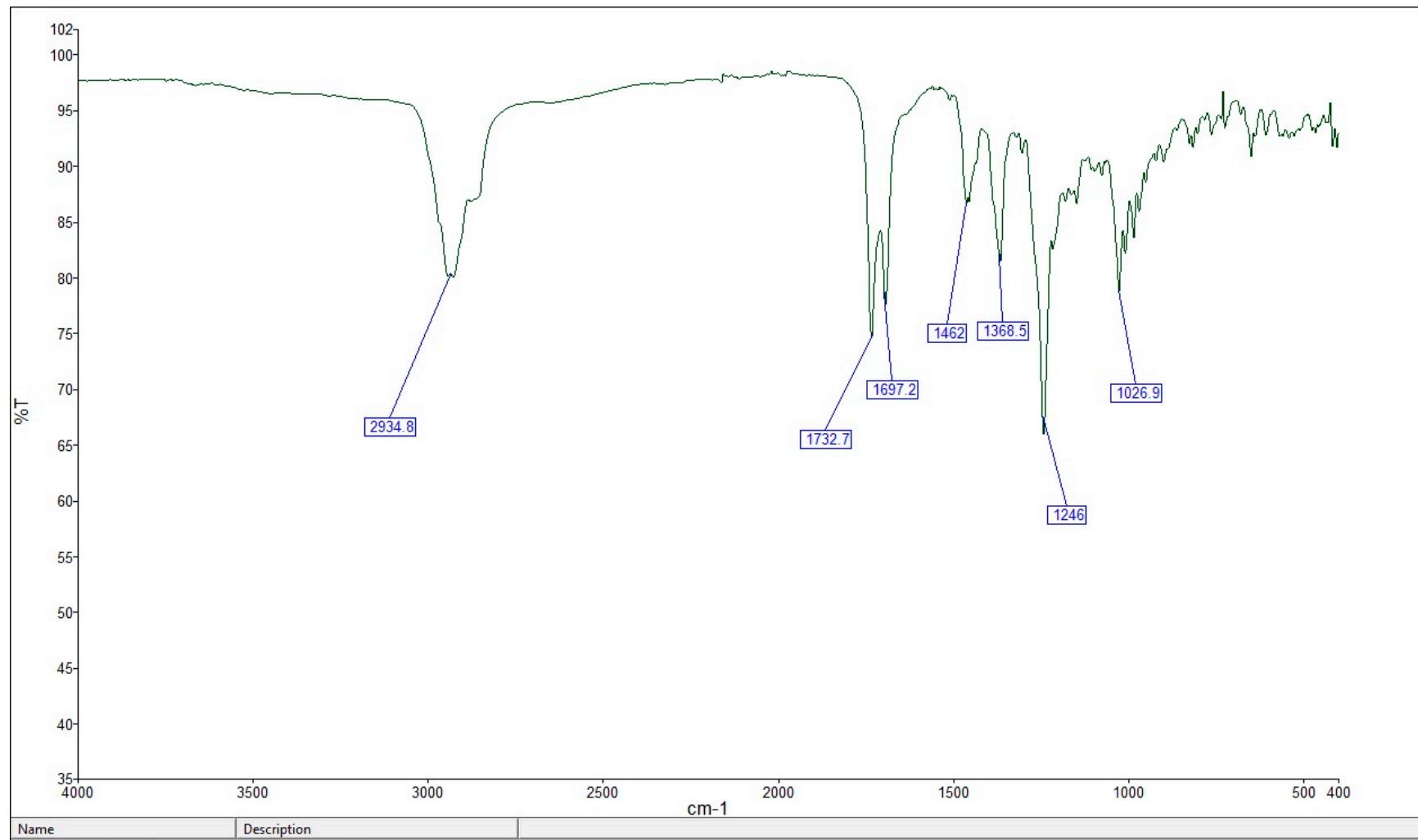


Figure S28: FT-IR spectrum of Oleanolic acid acetate (**7**).

PROTON_06
ZTPA266-274_2_new_cult

ppm	Intensity
1	7.63 408.4
2	7.59 434.9
3	7.26 704.6
4	7.08 188.4
5	7.08 232.6
6	7.06 228.7
7	7.06 323.5
8	7.04 501.7
9	7.03 387.8
10	6.92 642.5
11	6.90 505.2
12	6.31 619.0
13	6.27 566.7
14	4.20 382.8
15	4.19 839.7
16	4.17 398.8
17	3.93 3851.3
18	1.81 418.1
19	1.71 173.6
20	1.69 189.3
21	1.69 117.4
22	1.68 69.7
23	1.67 186.6
24	1.35 25.3
25	1.35 74.4
26	1.34 8.6
27	1.33 67.2
28	1.32 36.1
29	1.31 116.7
30	1.30 108.8
31	1.28 51.7
32	1.27 607.0
33	1.25 22579.1
34	1.17 276.9
35	1.09 263.8
36	0.89 324.2
37	0.88 1258.3
38	0.86 508.0

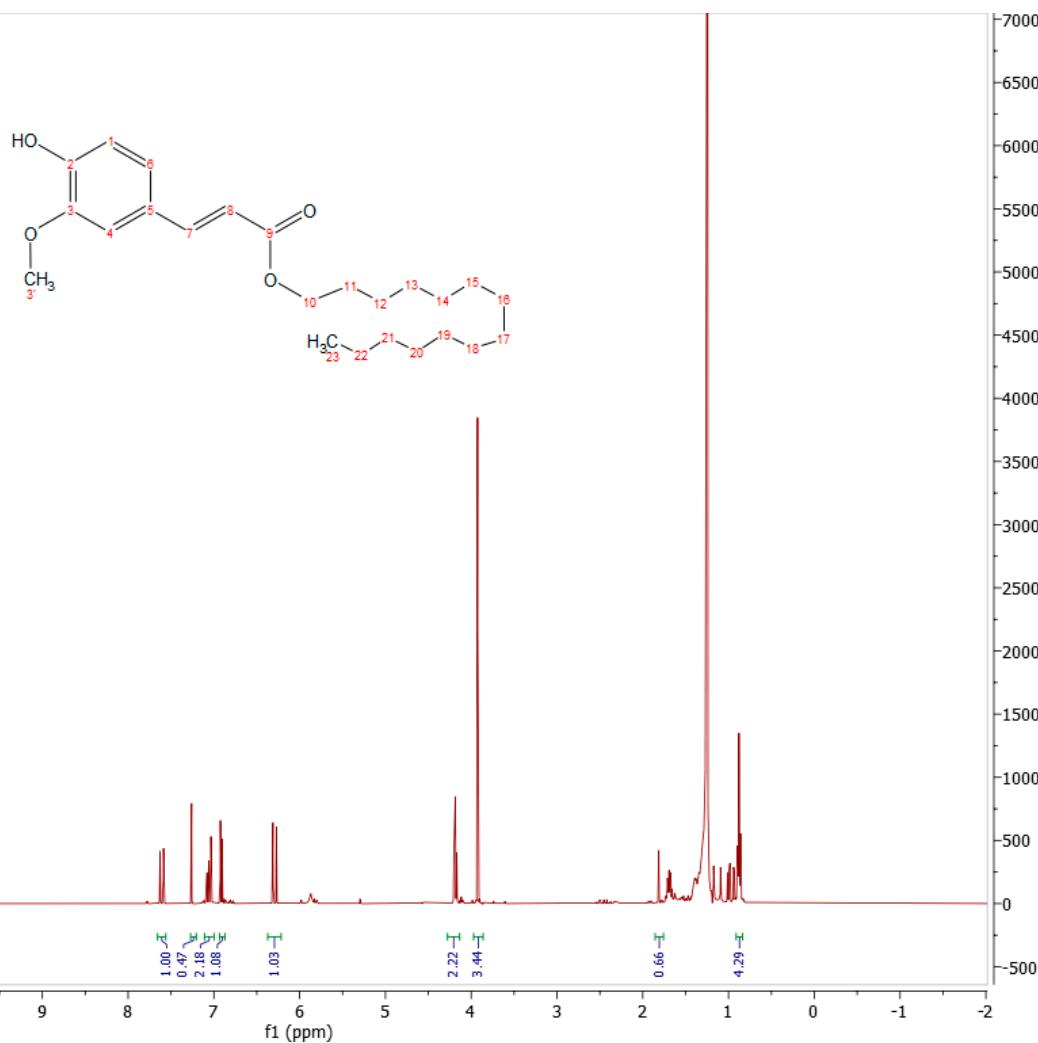


Figure S29: ¹H NMR spectrum of Tetradecyl (E)-ferulate (8) recorded in CDCl_3 .

CARBON_04
ZTPA266-274_2_new_cult

ppm	Intensity
1	167.3980 29.8
2	147.8672 29.4
3	146.7248 28.3
4	144.6238 53.1
5	127.0363 35.5
6	123.0440 74.2
7	115.6550 52.0
8	114.6752 66.6
9	109.2332 63.0
10	77.3329 171.3
11	77.0150 175.2
12	76.6983 166.9
13	64.6226 62.7
14	55.9126 70.4
15	31.9291 92.5
16	29.9972 6.9
17	29.7071 1140.1
18	29.6667 151.5
19	29.6095 75.1
20	29.5569 77.3
21	29.3708 92.4
22	29.3140 77.4
23	28.7676 69.0
24	28.1414 8.1
25	26.0005 75.6
26	22.6990 84.2
27	14.1320 72.7

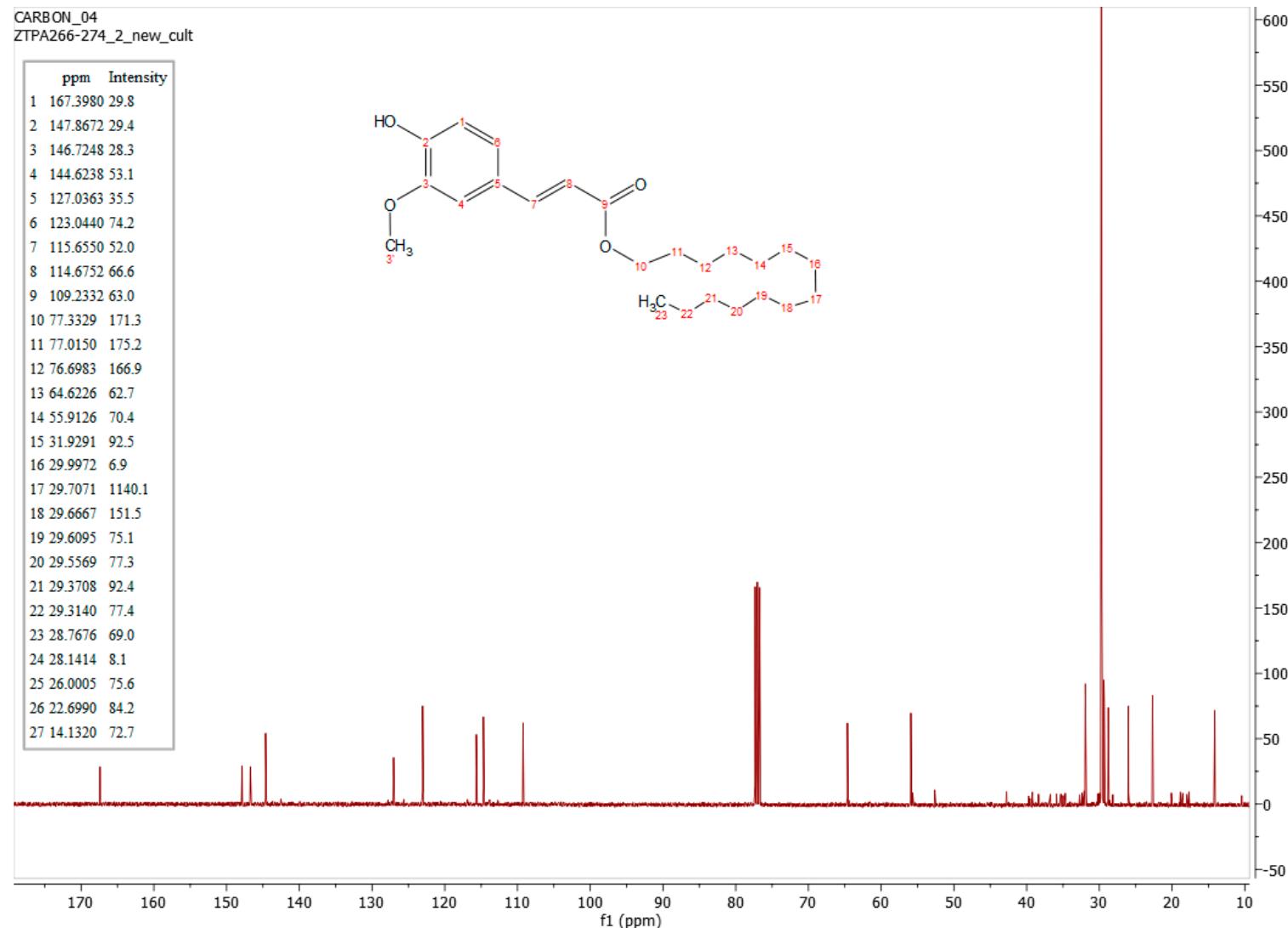
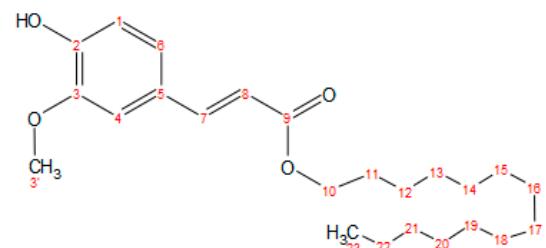


Figure S30: ^{13}C NMR spectrum of Tetradecyl (E)-ferulate (**8**) recorded in CDCl_3

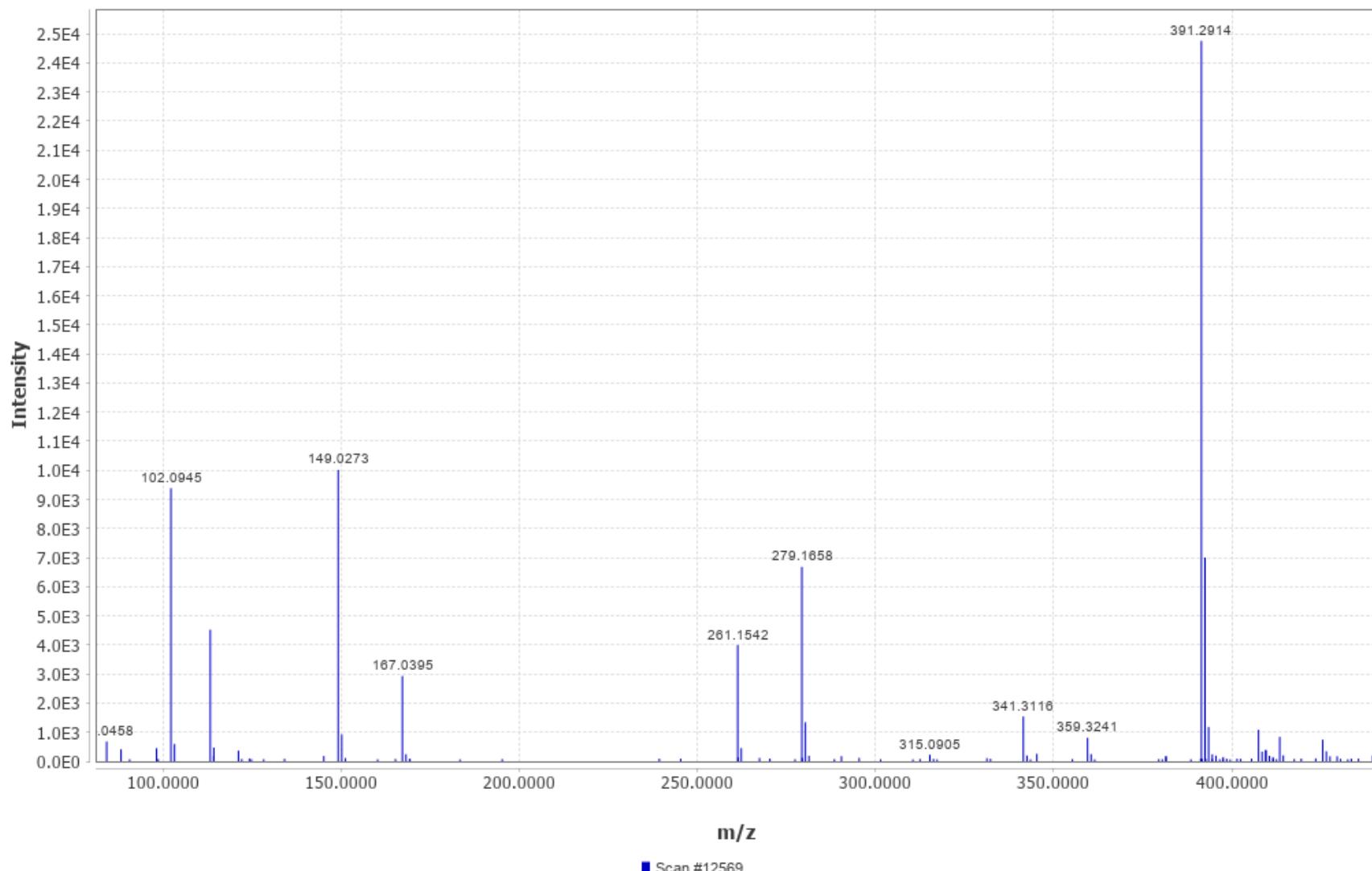


Figure S31: MS spectrum of Tetradecyl (E)-ferulate (**8**).

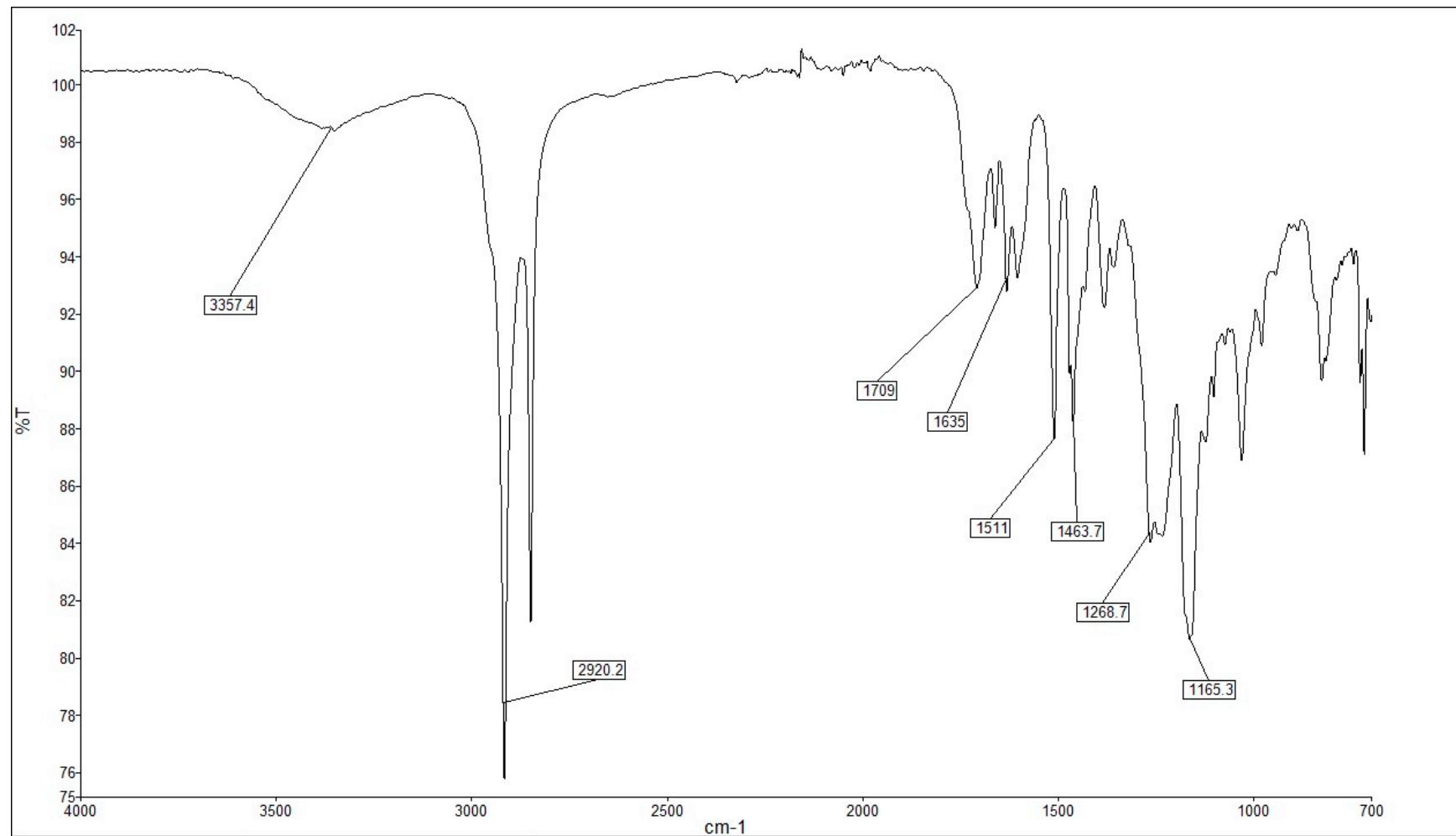


Figure S32: FT-IR spectrum of Tetradecyl (E)-ferulate (**8**).

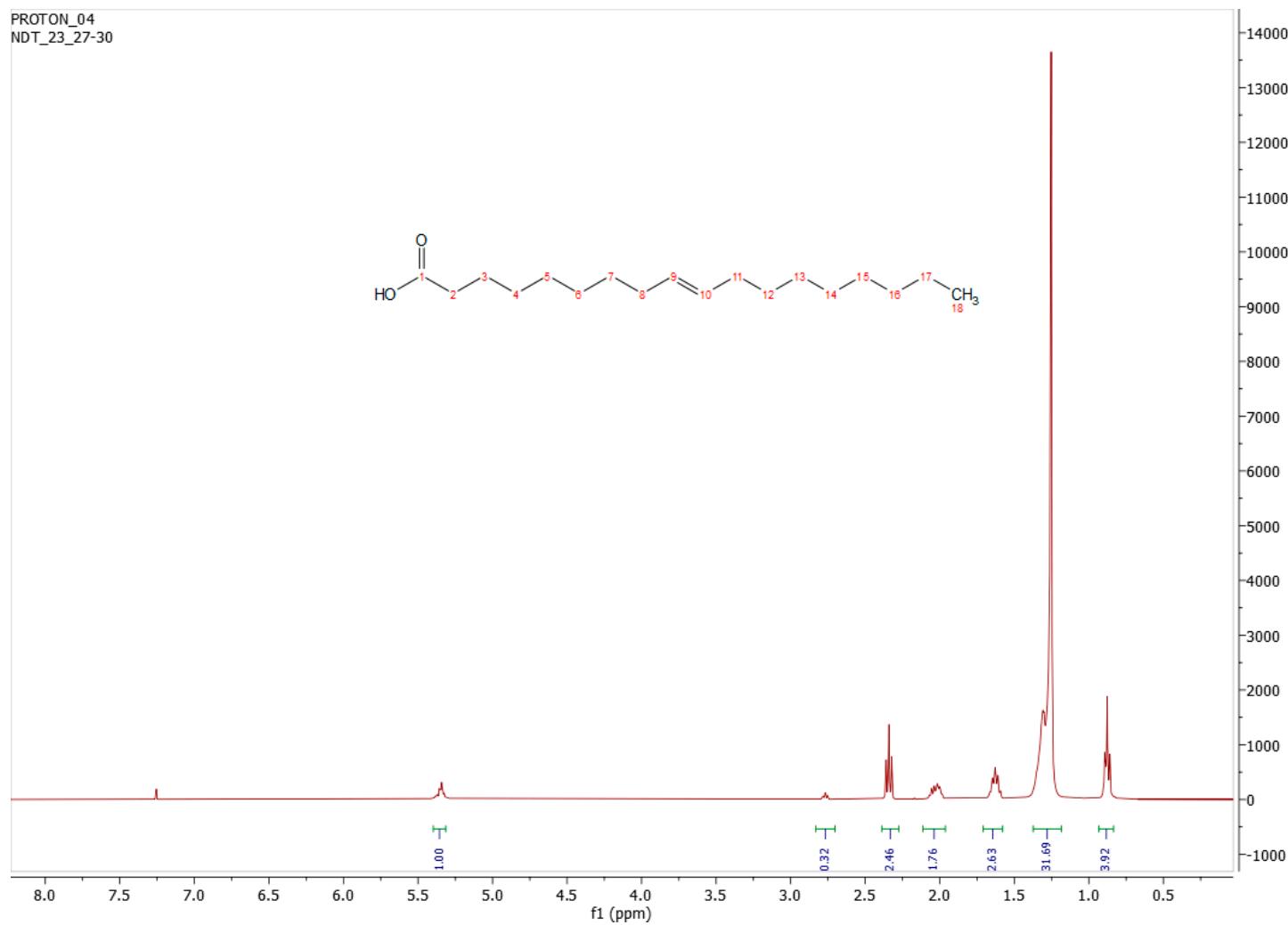


Figure S33: ^1H NMR spectrum of 9-Octadecenoic acid (**9**) recorded in CDCl_3 .

CARBON_03

	ppm	Intensity
1	180.1920	84.4
2	129.9973	135.0
3	128.0525	58.9
4	34.0496	264.9
5	31.9093	383.3
6	29.6826	1104.9
7	29.6477	494.5
8	29.6205	283.2
9	29.5710	396.2
10	29.4128	370.9
11	29.3431	361.5
12	29.3024	195.1
13	29.2190	394.6
14	29.1174	196.2
15	29.0444	516.4
16	24.6564	367.1
17	22.6689	401.1
18	14.0821	382.1

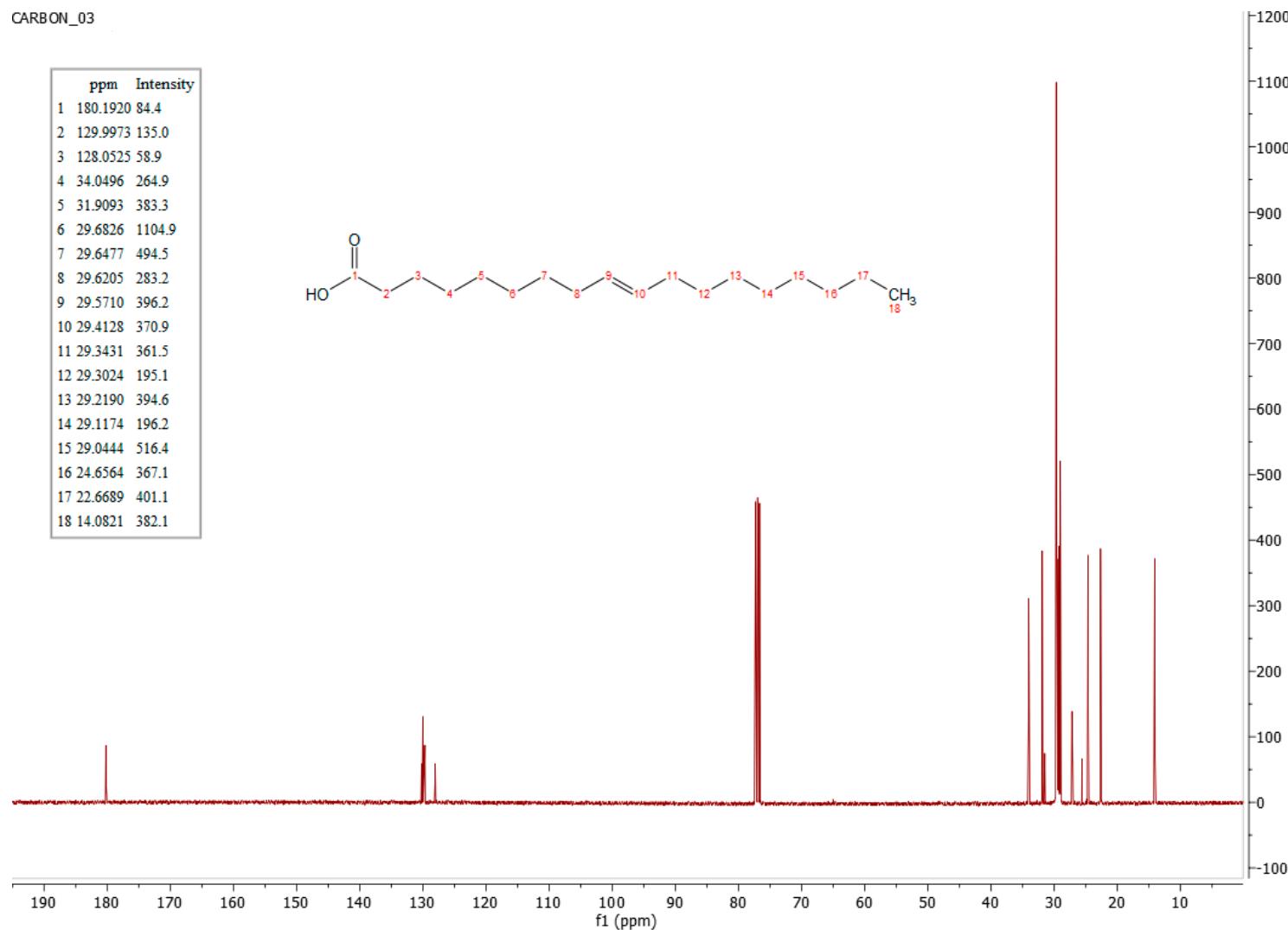
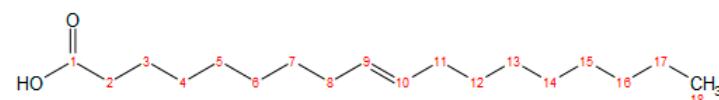


Figure S34: ^{13}C NMR spectrum of 9-Octadecenoic acid (**9**) recorded in CDCl_3

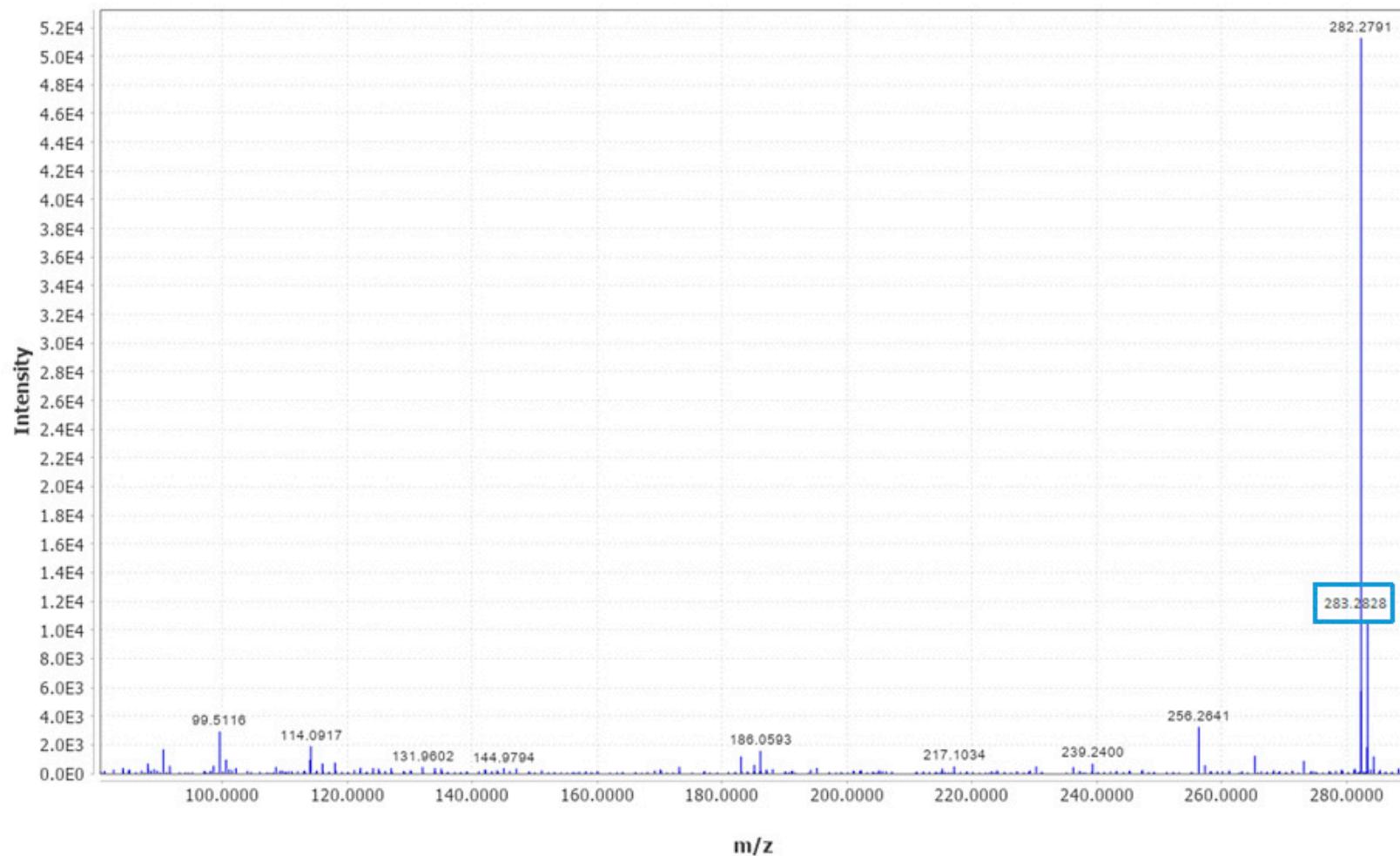


Figure S35: MS spectrum of 9-Octadecenoic acid (**9**).

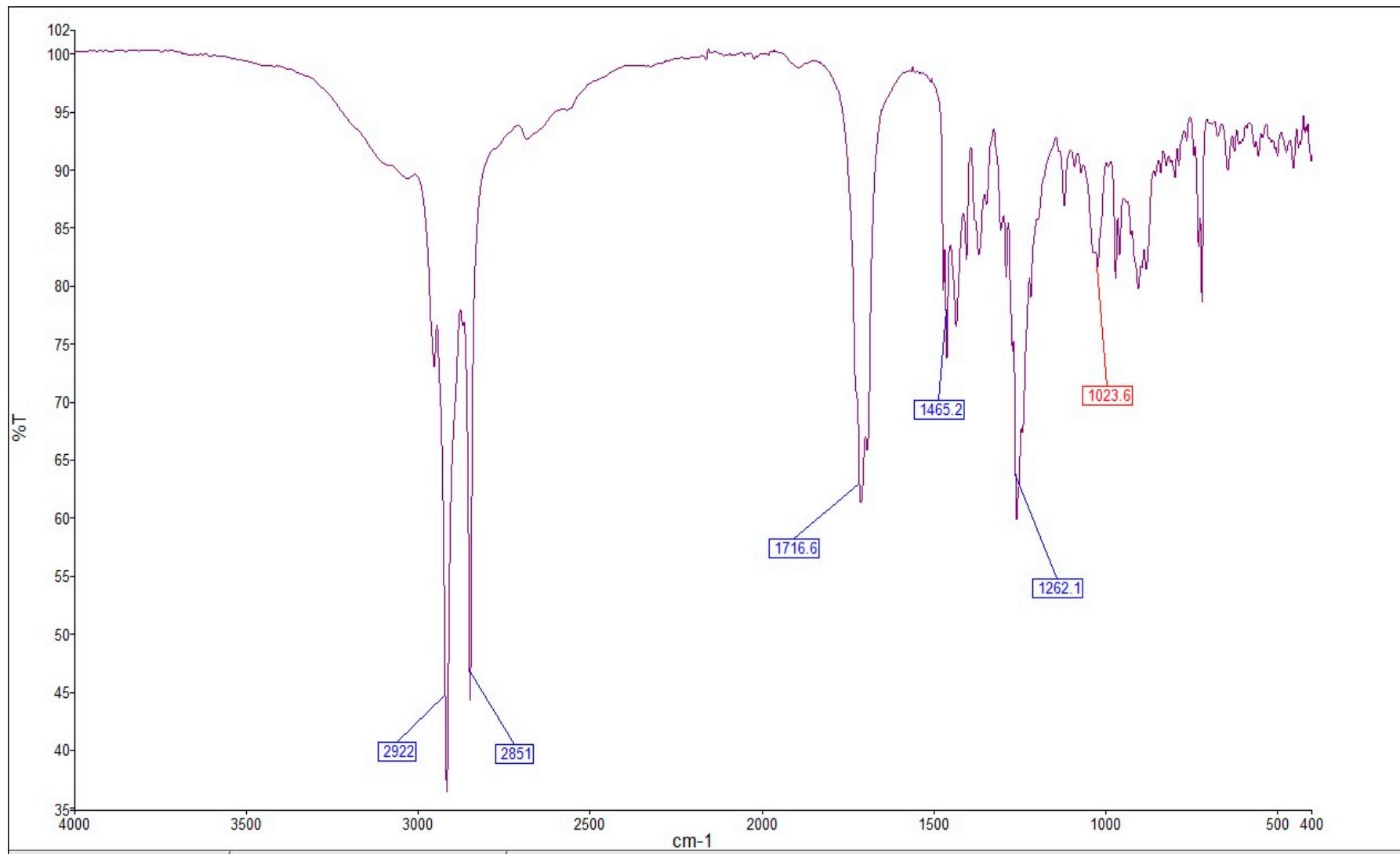


Figure S36: FT-IR spectrum of 9-Octadecenoic acid (**9**).

Table S1. ^{13}C NMR (100 MHz) data (δ value) of isolated compounds (1-9) in CHCl_3

C. No	13C NMR (100 MHz) δ value									
	Compound 1	Compound 2	Compound 3	Compound 4	Compound 5	Compound 6	Compound 7	Compound 8	Compound 9	
1.	22.3 (CH ₂)	36.1 (CH ₂)	32.1 (CH ₂)	38.7 (CH ₂)	37.2 (CH ₂)	113.0 (C)	38.1 (CH ₂)	127.0 (C)	180.1 (C)	
2.	41.6 (CH ₂)	211.9 (C)	195.0 (C)	27.4 (CH ₂)	31.7 (CH ₂)	165.9 (CH)	27.7 (CH ₂)	123.1 (CH)	34.2 (CH ₂)	
3.	213.1 (C)	76.9 (CH ₂)	142.6 (CH)	79.0 (CH)	71.8 (CH)	103.6 (CH)	80.9 (CH)	114.7 (CH)	24.7 (CH ₂)	
4.	58.2 (CH)	54.6 (CH)	140.8 (C)	38.9 (C)	42.3 (CH ₂)	55.2 (OCH ₃)	37.7 (C)	147.9(C)	28.9 (CH ₂)	
5.	42.1 (C)	38.1 (C)	39.7 (C)	55.3 (CH)	140.8 (C)	162.4 (C)	55.3 (CH)	146.7 (C)	29.3 (CH ₂)	
6.	41.6 (CH ₂)	40.6 (CH ₂)	38.4 (CH ₂)	18.3 (CH ₂)	121.7 (CH)	107.7 (CH)	18.2 (CH ₂)	109.3 (CH)	29.7 (CH ₂)	
7.	18.3 (CH ₂)	17.6 (CH ₂)	18.0 (CH ₂)	34.3 (CH ₂)	31.9 (CH ₂)	132.7 (CH)	32.5 (CH ₂)	144.8 (CH)	29.4 (CH ₂)	
8.	53.1 (CH)	53.9 (CH)	52.6 (CH)	40.8 (C)	31.9 (CH)	205.0 (C)	39.3 (C)	115.7 (CH)	27.2 (CH ₂)	
9.	37.4 (C)	37.6 (C)	36.7 (C)	50.4 (CH)	50.1 (CH)	46.0 (CH)	47.5 (CH)	167.4 (C)	130.0 (C)	
10.	59.5 (CH)	60.4 (CH)	55.7 (CH)	37.1 (C)	36.5 (C)	19.2 (CH ₃)	37.0 (C)	64 (CH ₂)	128.1 (C)	
11.	35.6 (CH ₂)	35.0 (CH ₂)	34.7 (CH ₂)	20.9 (CH ₂)	21.1 (CH)	133.5 (C)	23.4 (CH ₂)	28.9 (CH ₂)	25.7 (CH ₂)	
12.	30.5 (CH ₂)	30.1 (CH)	30.2 (CH ₂)	25.1 (CH ₂)	39.7 (CH ₂)	128.6 (CH)	122.6 (CH)	26.1 (CH ₂)	29.2 (CH ₂)	
13.	39.7 (C)	39.6 (C)	39.5 (C)	38.0 (CH)	42.2	114.4 (CH)	143.6 (C)	29.3 (CH ₂)	29.3 (CH ₂)	
14.	38.3 (C)	38.3 (C)	38.2 (C)	42.8 (C)	56.8 (CH)	55.2(OCH ₃)	41.6 (C)	29.4 (CH ₂)	29.6 (CH ₂)	
15.	32.4 (CH ₂)	32.2 (CH ₂)	32.2 (CH ₂)	27.4 (CH ₂)	24.4 (CH ₂)	158.7 (C)	27.7 (CH ₂)	29.5 (CH ₂)	29.7 (CH ₂)	
16.	36.0 (CH ₂)	35.8 (CH ₂)	35.9 (CH ₂)	35.6 (CH ₂)	29.7 (CH ₂)	114.4 (CH)	22.9 (CH ₂)	29.5 (CH ₂)	31.9 (CH ₂)	
17.	30.0 (C)	30.0 (C)	30.0 (C)	43.0 (C)	56.8 (CH)	128.6 (CH)	46.5 (C)	29.6 (CH ₂)	22.7 (CH ₂)	
18.	42.8 (CH)	42.7 (CH)	42.6 (CH)	48.3 (CH)	12.0 (CH ₃)	-	41.0 (CH ₂)	29.6 (CH ₂)	14.1 (CH ₃)	
19.	35.3 (CH)	35.3 (CH ₂)	35.3 (CH ₂)	48.0 (CH)	19.4 (CH ₃)	-	45.8 (CH ₂)	29.7 (CH ₂)	-	
20.	28.1 (C)	28.1 (CH ₂)	28.1 (C)	151.0 (C)	40.5 (CH)	-	30.7 (C)	29.7 (CH ₂)	-	
21.	32.8 (CH ₂)	32.8 (CH ₂)	32.7 (CH ₂)	29.8 (CH ₂)	21.2 (CH ₃)	-	33.8 (CH ₂)	29.8 (CH ₂)	-	
22.	39.2 (CH ₂)	39.2 (CH ₂)	39.3 (CH ₂)	40.0 (CH ₂)	138.3 (CH)	-	32.4 (CH ₂)	22.7 (CH ₂)	-	
23.	6.8 (CH ₃)	10.8 (CH ₃)	10.4 (CH ₃)	28.0 (CH ₃)	129.3 (CH)	-	16.6 (CH ₃)	14.2 (CH ₃)	-	

24.	14.6 (CH ₃)	14.2 (CH ₃)	17.7 (CH ₃)	15.3 (CH ₃)	51.2 (CH)	-	28.0 (CH ₃)	55.9 (OCH ₃)	-
25.	17.9 (CH ₃)	17.6 (CH ₃)	18.6 (CH ₃)	16.1 (CH ₃)	31.9 (CH)	-	15.4 (CH ₃)	-	-
26.	20.2 (CH ₃)	19.9 (CH ₃)	18.9 (CH ₃)	16.0 (CH ₃)	21.1 (CH ₃)	-	17.1 (CH ₃)	-	-
27.	18.6 (CH ₃)	18.6 (CH ₃)	20.1 (CH ₃)	14.5 (CH ₃)	19.0 (CH ₃)	-	25.9 (CH ₃)	-	-
28.	32.1 (CH ₃)	32.2 (CH ₃)	32.1 (CH ₃)	18.0 (CH ₃)	25.4 (CH ₂)	-	183.4 (C)	-	-
29.				109.3 (CH ₂)		-	23.6 (CH ₃)	-	-
	35.0 (CH ₃)	31.9 (CH ₃)	31.8 (CH ₃)		12.2 (CH ₃)				
30.	31.8 (CH ₃)	35.3 (CH ₃)	35.0 (CH ₃)	19.3 (CH ₃)	-	-	33.0 (CH ₃)	-	-
31.	-	-	-	-	-	-	21.3 (CH ₃)	-	-
32.	-	-	-	-	-	-	171.0 (C)	-	-

Table S2. ^1H NMR (400 MHz) data (δ value) of isolated compounds (1-5 and 7) in CHCl_3

C. No	^1H NMR (400MHz, δ , ppm)					
	Compound 1	Compound 2	Compound 3	Compound 4	Compound 5	Compound 7
1.	1.91 (1H, m) 1.65 (1H, m)	2.39 (1H, t, $J=13.8$ Hz) 2.52 (1H, dd $J= 10.8, 2.9$ Hz)	2.42 (1H, dd $J= 17.8.0, 15.7$ Hz) 2.53 (1H, dd $J= 17.8, 3.64$ Hz)	0.91 (1H, m)	1.85 (1H, m), 1.08 (1H, m)	1.60, 1.05 (2H, m)
2.	2.36 (1H, m) 2.27 (1H, m)	-	-	1.67 (1H, m)	1.84 (1H, m), 1.50 (1H, m)	162, 188 (2H, m)
3.	-	3.81 (1H, dd $J=11.8, 2.96$ Hz) 3.51 (1H, d, $J= 3.48$ Hz OH-3)	5.98 (1H, s, OH-3)	1.60 (1H, d, $J = 10.0$ Hz)	3.51 (tdd, $J = 6.1, 4.4, 5.1$ Hz)	4.51 (1H, t, $J= 7.6$ Hz)
4.	2.21 (1H, q, $J=6.7$ Hz)	1.29 (m)	-	1.52 (1H, m)	2.30 (1H, m) 2.23 (1H, m)	-
5.	-	-	-	3.19 (1H, dd, $J = 4.4, 6.4$ Hz)	-	0.85 (1H, m)
6.	1.73 (1H, m) 1.23 (1H, m)	1.06 (1H, m) 1.85 (1H, br d $J=12.8$ Hz)	1.93 (1H, dd $J= 12.0, 2.52$ Hz) 181 (1H, m)	-	5.34 (d, $J = 5.2$ Hz)	1.52, 1.40 (2H, m)
7.	1.47 (1H, m) 1.38 (1H, m)	1.43 (1H, m) 1.48 (1H, m)	1.50 (1H, m) 1.44 (1H, m)	0.70 (1H, d, $J = 5.2$ Hz)	1.99(1H, m) 1.53 (1H, m)	1.40, 1.28 (2H, m)
8.	1.38 (1H, m)	1.29 (1H, m)	1.37 (1H, m)	1.52 (1H, m)	1.49 (1H, m)	-
9.	-	-	-	1.39 (1H, m)	0.93 (1H, m)	1.55 (1H, t)
10.	1.53 (1H, m)	1.30 (1H, m)	1.80 (1H, m)	<1.38> (2H, m)	-	-
11.	1.45 (1H, m) 1.27 (1H, m)	1.20 (1H, m) 1.30 (1H, m)	1.32 (1H, m) 1.26 (1H, m)	-	<1.50> (1H, m)	1.88, 1.62 (2H, dd, $J= 6.0, 3.7$ Hz)
12.	1.32 <1H, m>	1.33 (2H, m)	1.33 (2H, m)	1.26 (1H, m)	2.04 (1H, m) 1.17 (H, m)	5.28 (1H, t, $J= 3.7$ Hz)
13.	-	-	-	-	-	-

14.	-	-	-	1.40 (1H, m)	1.02 (1H, m)	-
15.	1.46 (1H, m) 1.25 (1H, m)	1.27 (1H, m) 1.51 (1H, m)	1.50 (1H, m) 1.30 (1H, m)	1.25 (1H, m)	1.56 (1H, m) 1.03 (1H, m)	1.05, 172 (2H, m)
16.	1.55 (1H, m)	1.33 (1H, m) 1.51 (1H, m)	1.54 (1H, m) 1.38 (1H, m)	1.67 (1H, d, <i>J</i> =9.6Hz),	1.26 (1H, m) 1.05 (1H, m)	1.98, 1.94 (2H, dd, <i>J</i> = 9.4, 4.0 Hz)
17.	-	-	-	1.02 (1H, m)	1.16 (1H, m)	-
18.	1.53 (1H, m)	1.52 (1H, m)	1.55 (1H, m)	1.67 (1H, t)	0.69 (3H, s)	2.84 (2H, dd, <i>J</i> = 9.4, 4.6 Hz)
19.	1.44 (1H, m)	1.16 (1H, m) 1.32 (1H, m)	1.20 (1H, m)	-	1.02 (3H, s)	1.62, 1.18 (2H, m)
20.	-	-	-	1.71 (1H, m)	2.04 (1H, m)	-
21.	1.46 (1H, m) 1.26 (1H, m)	1.47 (2H, m)	1.46 (1H, m) 1.28 (1H, m)	1.02 (1H, m)	1.02 (3H, s)	1.22, 1.19 (2H, m)
22.	1.45 (1H, m) 0.93 (1H, m)	1.46 (2H, m)	1.51 (1H, m)	1.46 (1H, m)	5.14 (1H, dd, <i>J</i> = 8.5, 6.6 Hz)	1.76, 1.57 (2H, m)
23.	0.86 (3H, d, <i>J</i> =6.3 Hz)	1.05 (3H, d, <i>J</i> =6.6 Hz)	1.80 (3H, s)	1.36 (1H, m)	5.05 (1H, dd, <i>J</i> = 8.5, 6.4 Hz)	0.85 (3H, s)
24.	0.71 (3H, s)	1.03 (3H, s)	0.93 (3H, s)	-	1.52 (1H, m)	0.86 (3H, s)
25.	0.86 (3H, s)	0.87 (3H, s)	0.98 (3H, s)	1.36 (1H, t)	1.52 (1H, m)	0.94 (3H, s)
26.	0.98 (3H, s)	1.00 (3H, s)	1.09 (3H, s)	2.38 (1H, td, <i>J</i> = 5.2,5.6 Hz)	0.83 (3H, d, <i>J</i> = 5.8 Hz)	0.75 (3H, s)
27.	1.04 (3H, s)	0.97 (3H, s)	1.00 (3H, s)	-	0.80 (3H, d, <i>J</i> = 5.8 Hz)	1.13 (3H, s)
28.	1.16 (3H, s)	1.17 (3H, s)	1.17 (3H, s)	1.91 (1H, m)	1.42 (1H, m) 1.16 (1H, m)	-
29.	0.99 (3H, s)	0.98 (3H, s)	0.99 (3H, s)	1.33 (1H, m)	0.80 (3H, t, <i>J</i> = 6.4 Hz)	0.93 (3H, s)
30.	0.94 (3H, s)	0.94 (3H, s)	0.93 (3H, s)	1.38 (1H, m)	-	0.90 (3H, s)
31.	-	-	-	-	-	2.04 (3H, s)

Table S3. ^1H NMR (400 MHz) data (δ value) of isolated compounds (6, 8 and 9) in CHCl_3

Carbon No	^1H NMR (400MHz, δ , ppm)		
	Compound 6	Compound 8	Compound 9
1.	-	-	-
2.	13.05 (1H, s, -OH)	7.07 (1H, dd, $J= 8.2, 1.6$ Hz)	2.34 (2H, t, $J= 7.5$ Hz)
3.	6.36 (1H, d, $J= 2.6$ Hz)	6.92 (1H, d, $J= 8.1$ Hz)	1.63 (2H, m)
4.	3.76 (3H, s, OCH_3)	-	1.31 (2H, m)
5.		-	1.26 (2H, m)
6.	6.33 (1H, dd, $J= 8.0, 2.6$ Hz)	7.03 (1H, d, $J= 2.0$ Hz)	1.26 (2H, m)
7.	7.70 (1H, d, $J= 8.0$ Hz)	7.36 (1H, d, $J= 15.9$ Hz)	1.30 (2H, m)
8.	-	6.31(1H, d, $J= 15.9$ Hz)	2.02 (2H, m)
9.	4.60 (1H, q, $J= 8.0$ Hz)	-	5.35 (1H, m)
10.	1.50 (3H, s)	4.19 (2H, t, $J= 6.7$ Hz)	5.35 (1H, m)
11.	-	1.69 (2H, m)	2.78 (2H, m)
12.	7.21 (1H, d, $J= 7.0$ Hz)	1.38 (2H, m)	1.25 (2H, m)
13.	6.85 (1H, d, $J= 7.0$ Hz)	1.34 (2H, m)	1.26 (2H, m)
14.	3.76 (3H, s, OCH_3)	1.33 (2H, m)	1.26 (2H, m)
15.	-	1.32 (2H, m)	1.26 (2H, m)
16.	6.83 (2H, d, $J= 7.0$ Hz)	1.31 (2H, m)	1.27 (2H, m)
17.	7.21 (2H, d, $J= 7.0$ Hz)	1.32 (2H, m)	1.30 (2H, m)
18.	-	1.30 (2H, m)	0.88 (3H, t, $J= 7.0$ Hz)
19.	-	1.26 (2H, m)	-
20.	-	1.26 (2H, m)	-
21.	-	1.25 (2H, m)	-
22.	-	1.25 (2H, m)	-
23.	-	3.91 O- CH_3 (3H, s)	-

24.	-	0.88 (3H, t, $J= 6.6$ Hz)	-
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