

Supporting Information

Figure S1. $^1\text{H-NMR}$ for loddigesinol G (1).

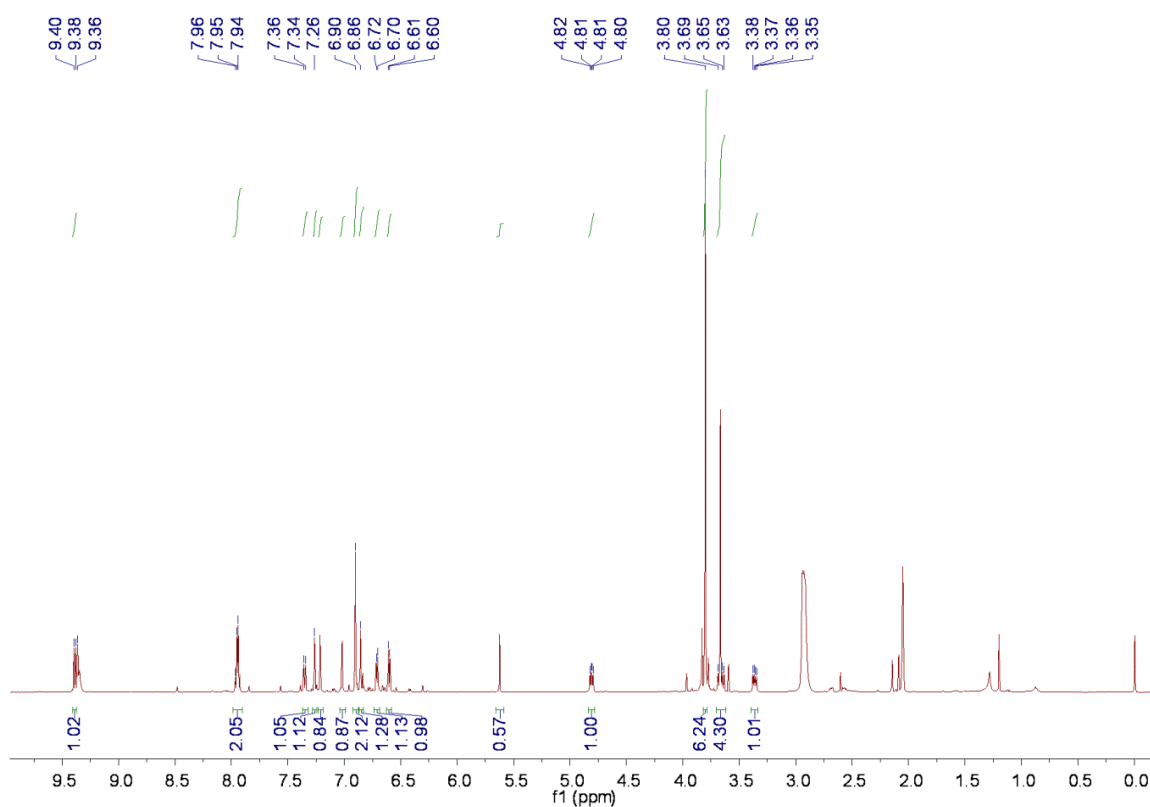


Figure S2. $^{13}\text{C-NMR}$ for loddigesinol G (1).

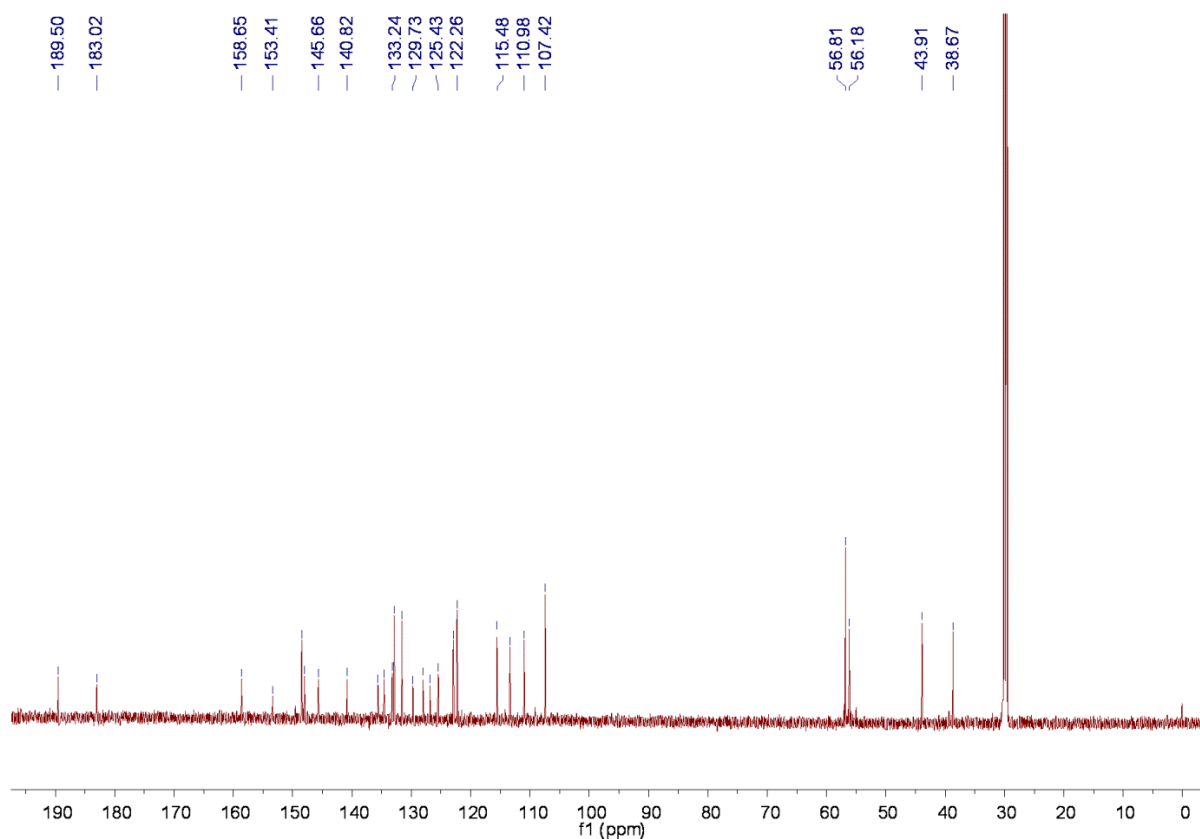


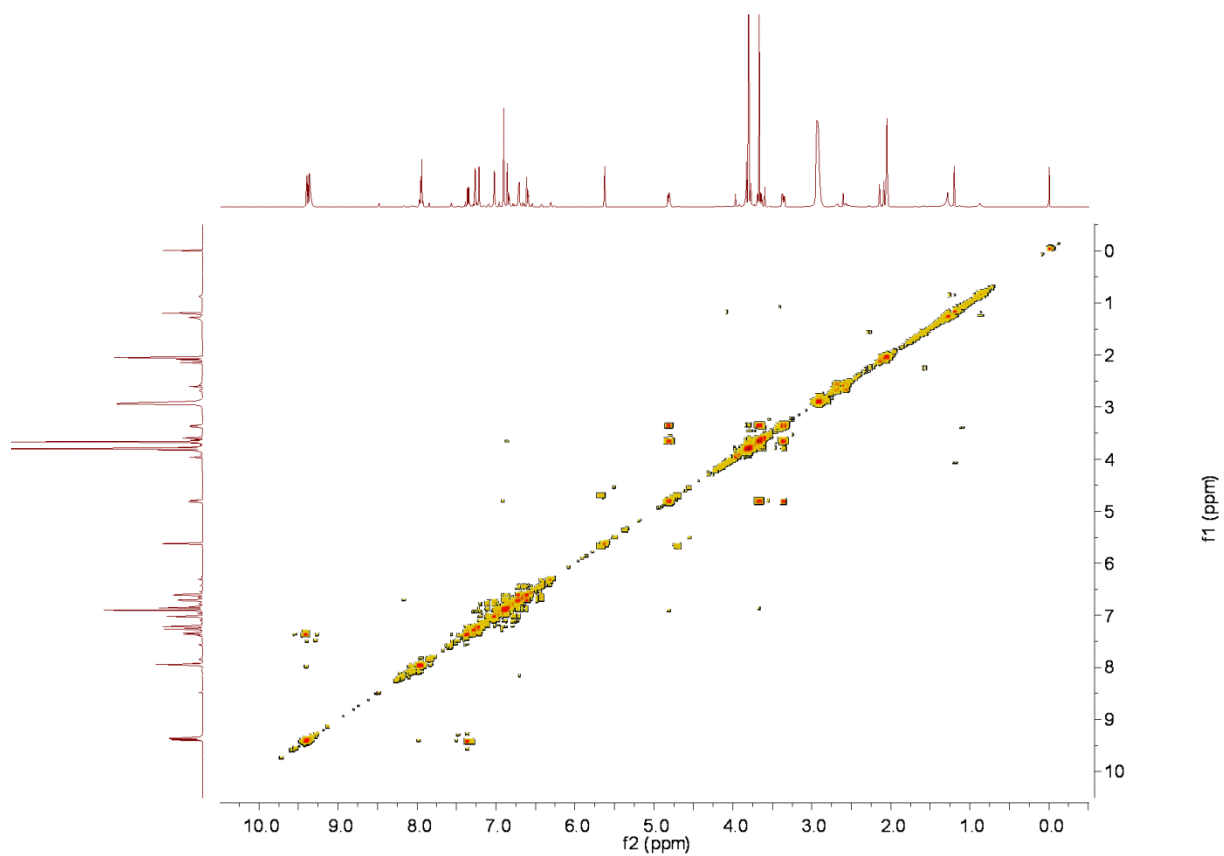
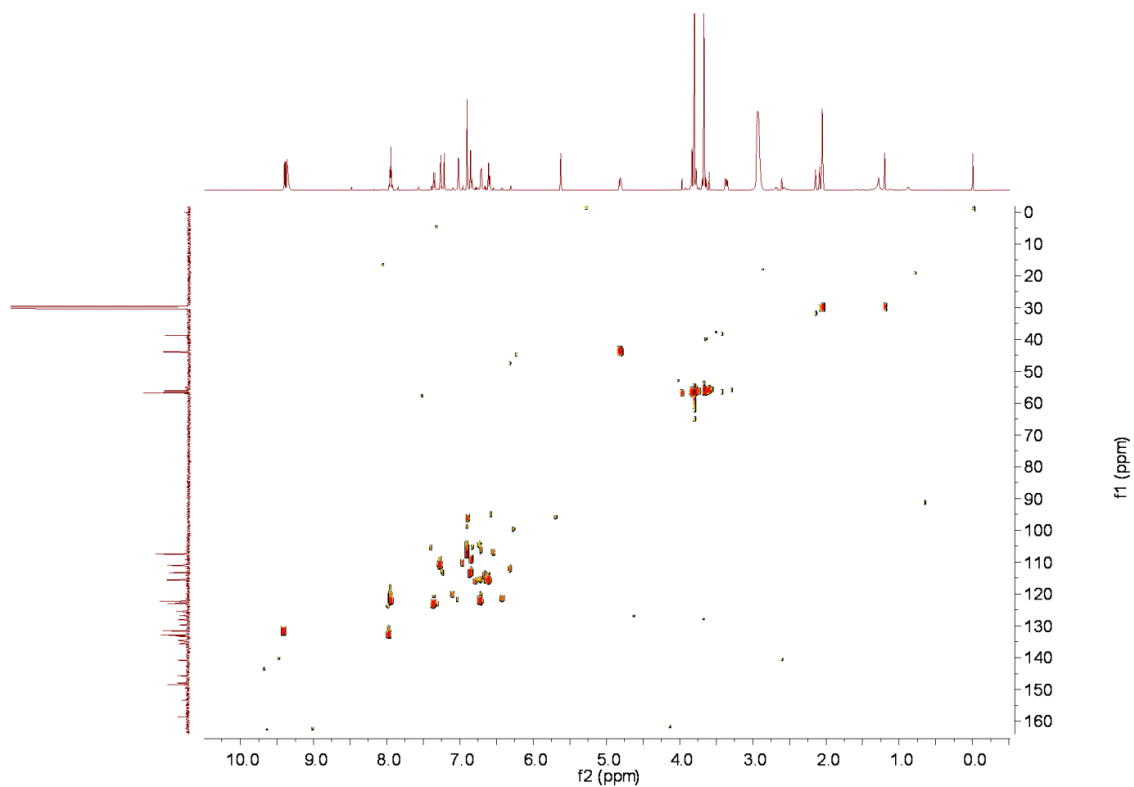
Figure S3. ^1H - ^1H COSY for loddigiinol G (**1**).**Figure S4.** HSQC for loddigiinol G (**1**).

Figure S5. HMBC for loddigesinol G (1).

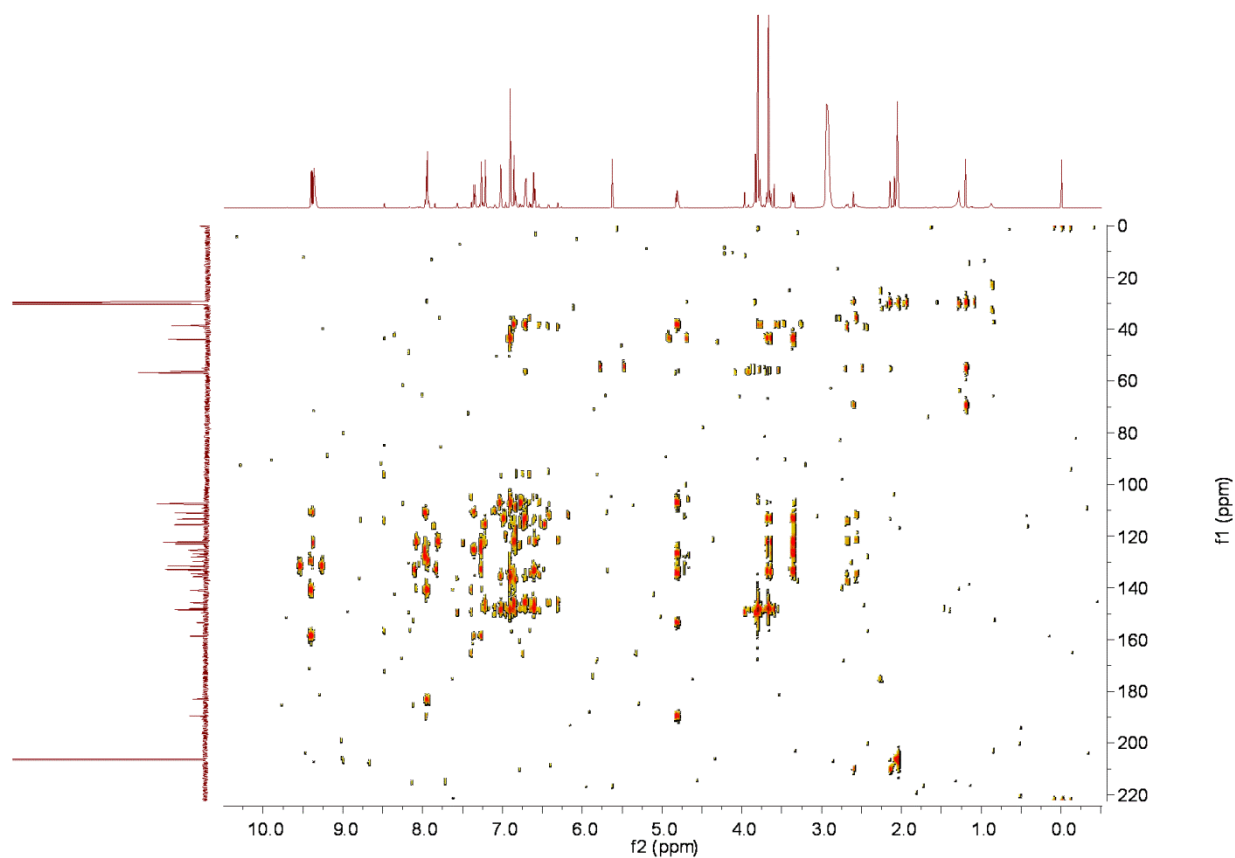
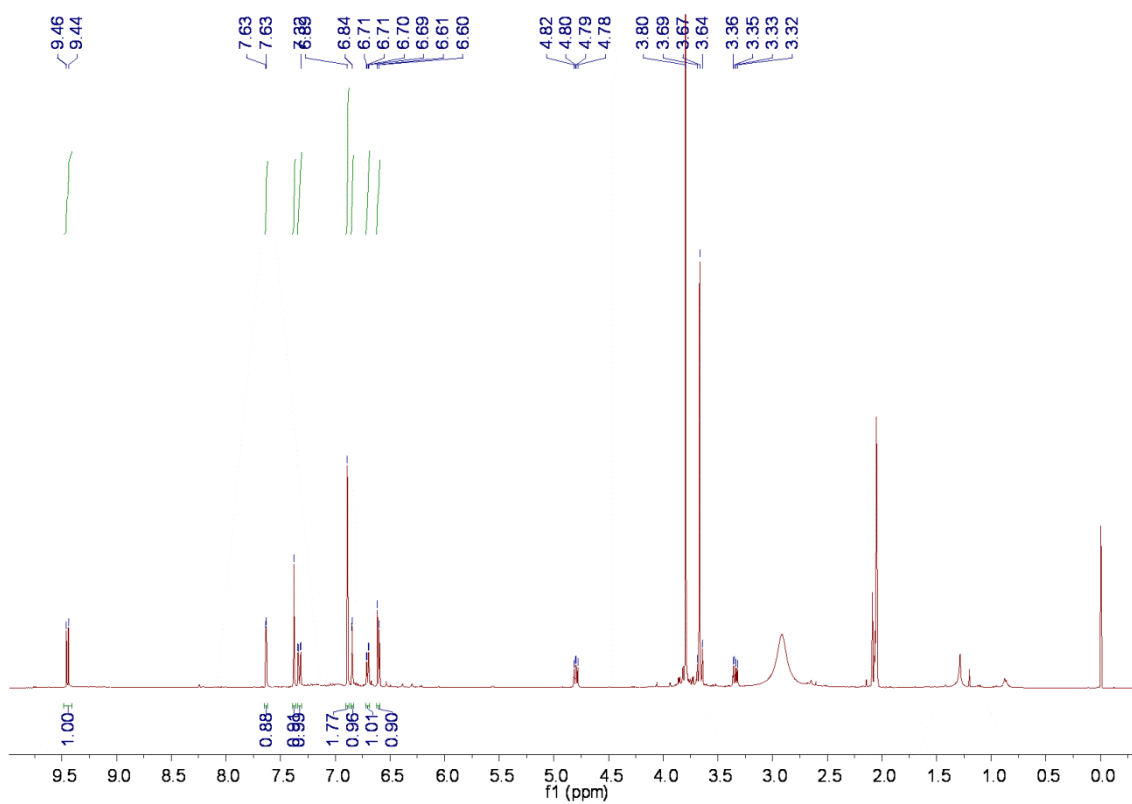
Figure S6. ^1H -NMR for loddigesinol H (2).

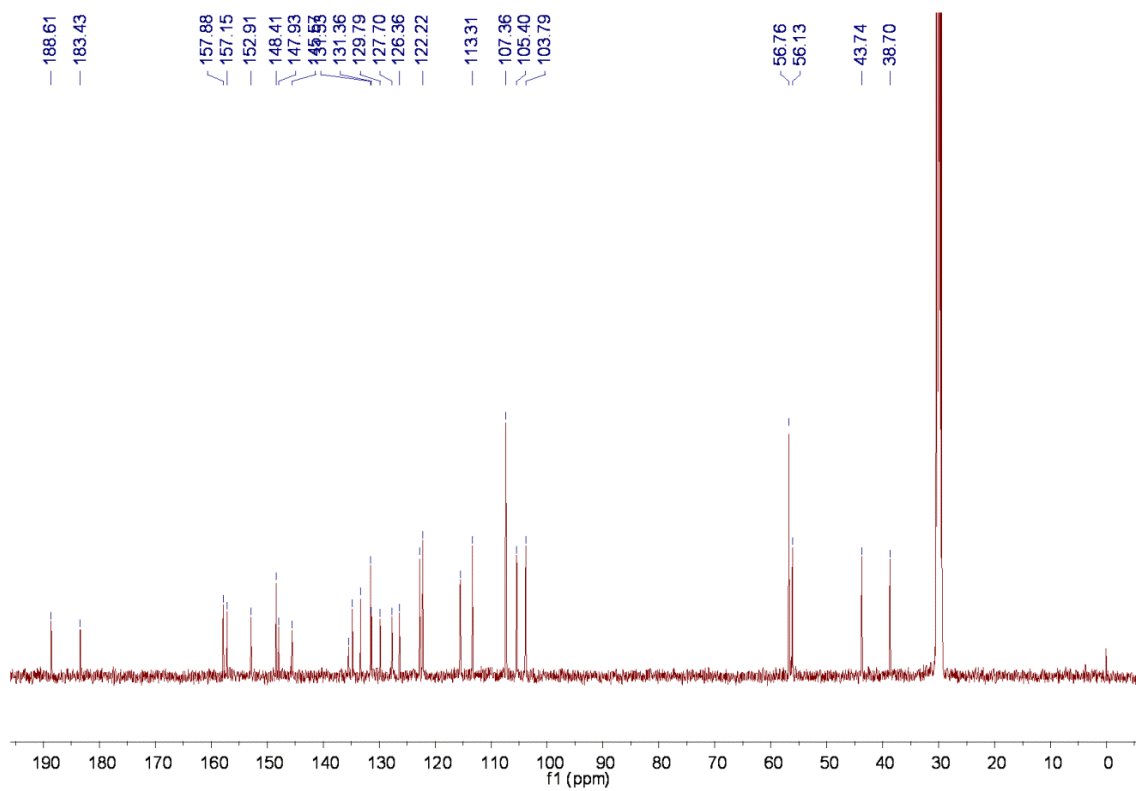
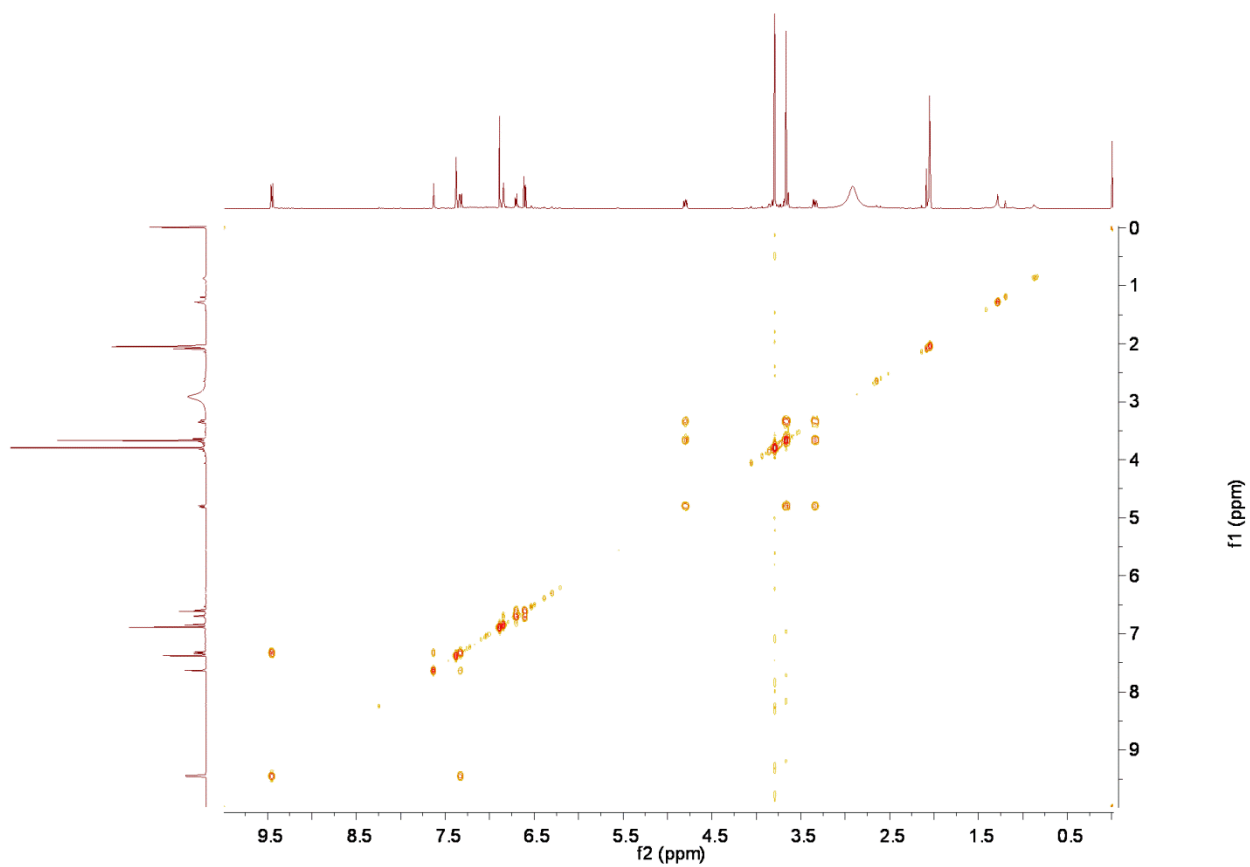
Figure S7. ^{13}C -NMR for loddigesinol H (2).**Figure S8.** ^1H - ^1H COSY for loddigesinol H (2).

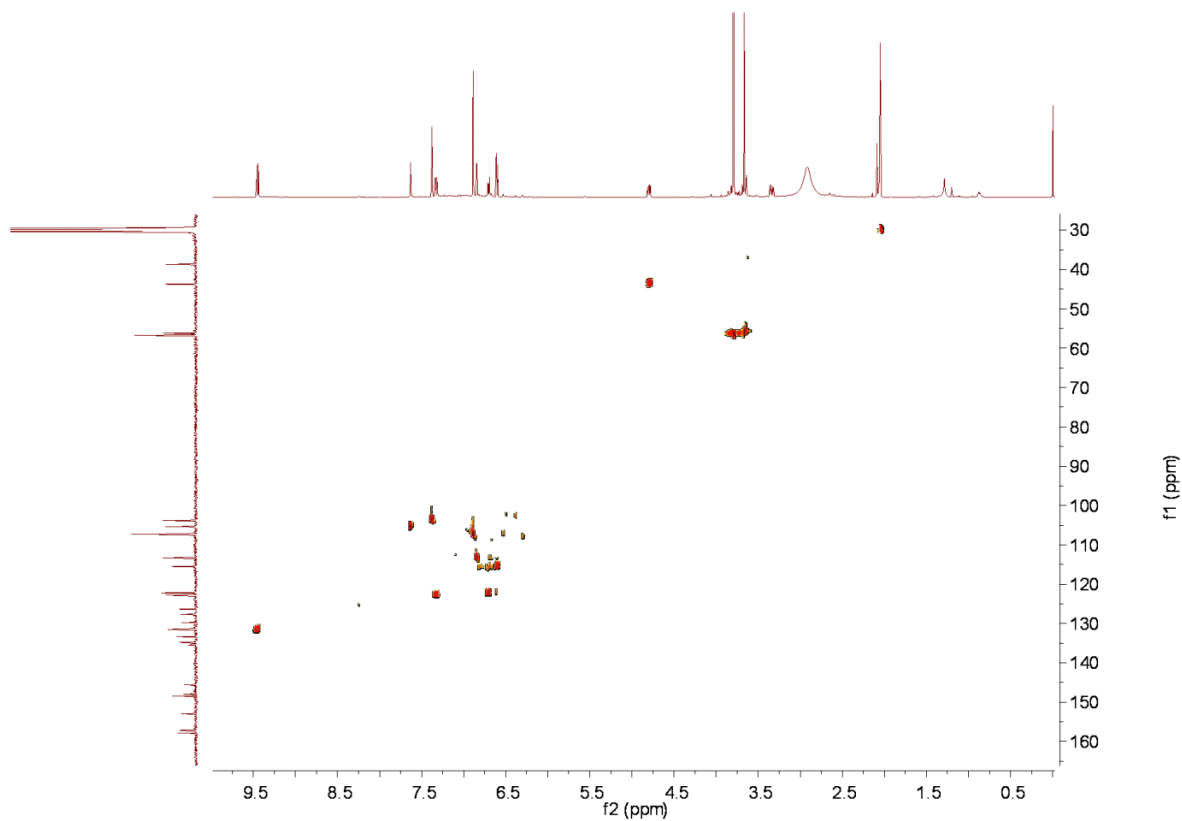
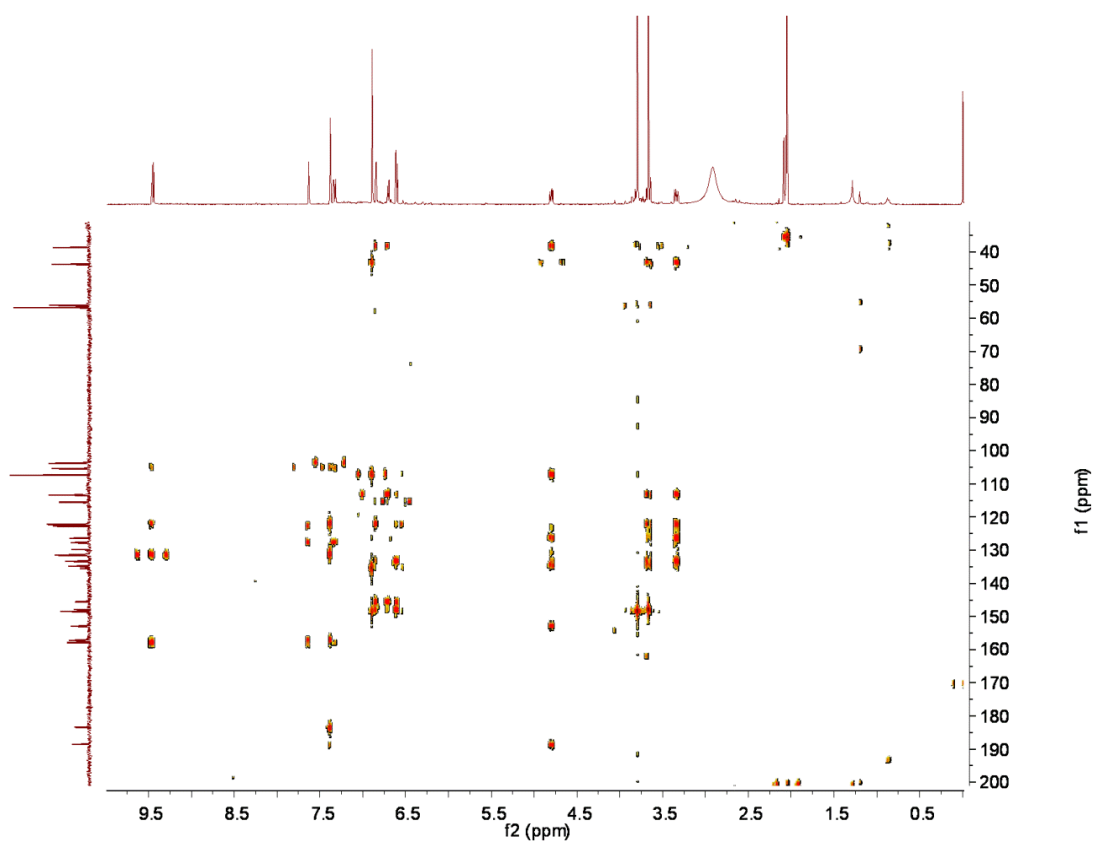
Figure S9. HSQC for loddigesiinol H (2).**Figure S10.** HMBC for loddigesiinol H (2).

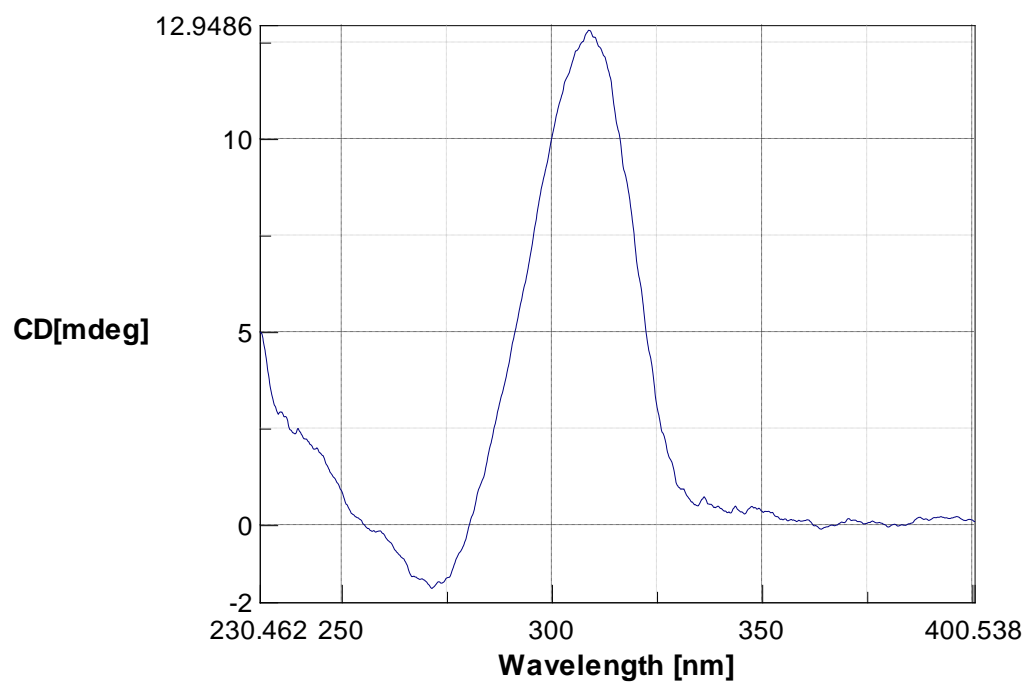
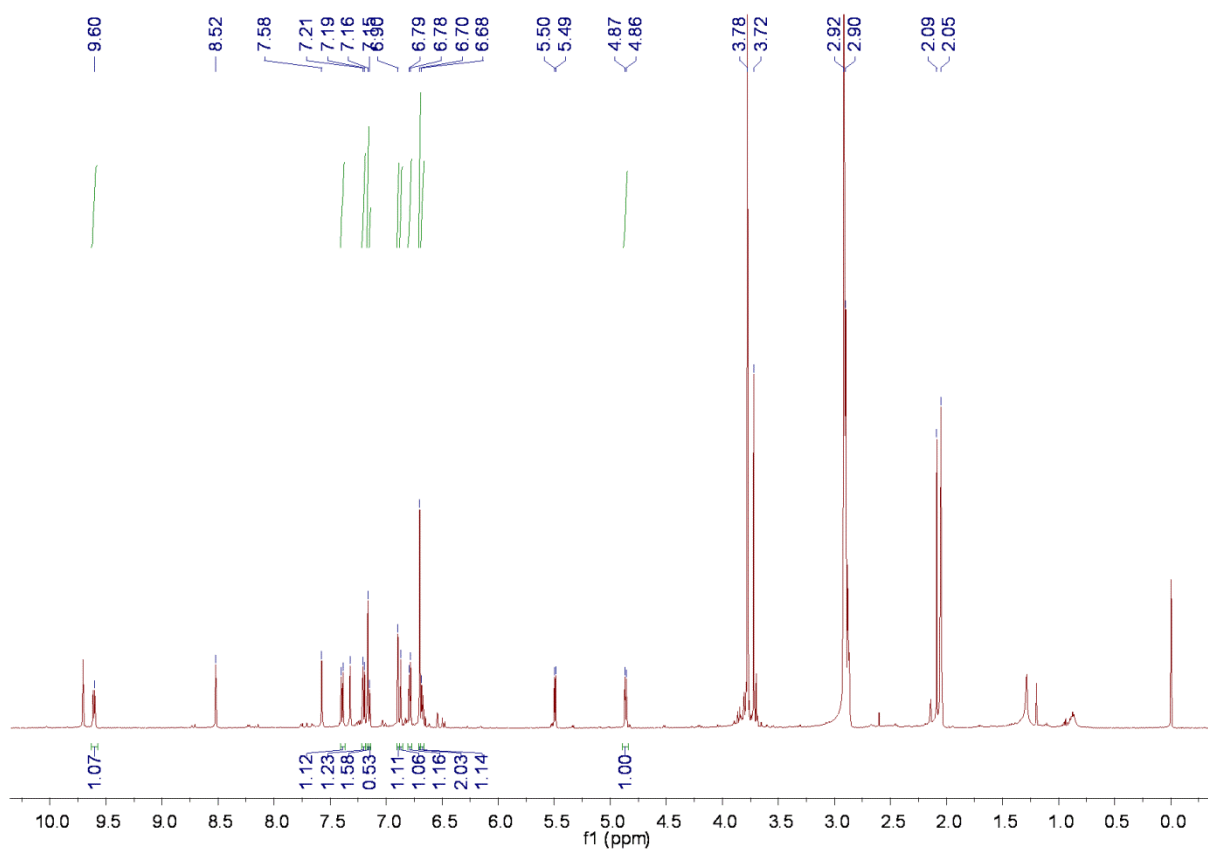
Figure S11. ECD spectrum for loddigesiinol H (2).**Figure S12.** $^1\text{H-NMR}$ for loddigesiinol I (3).

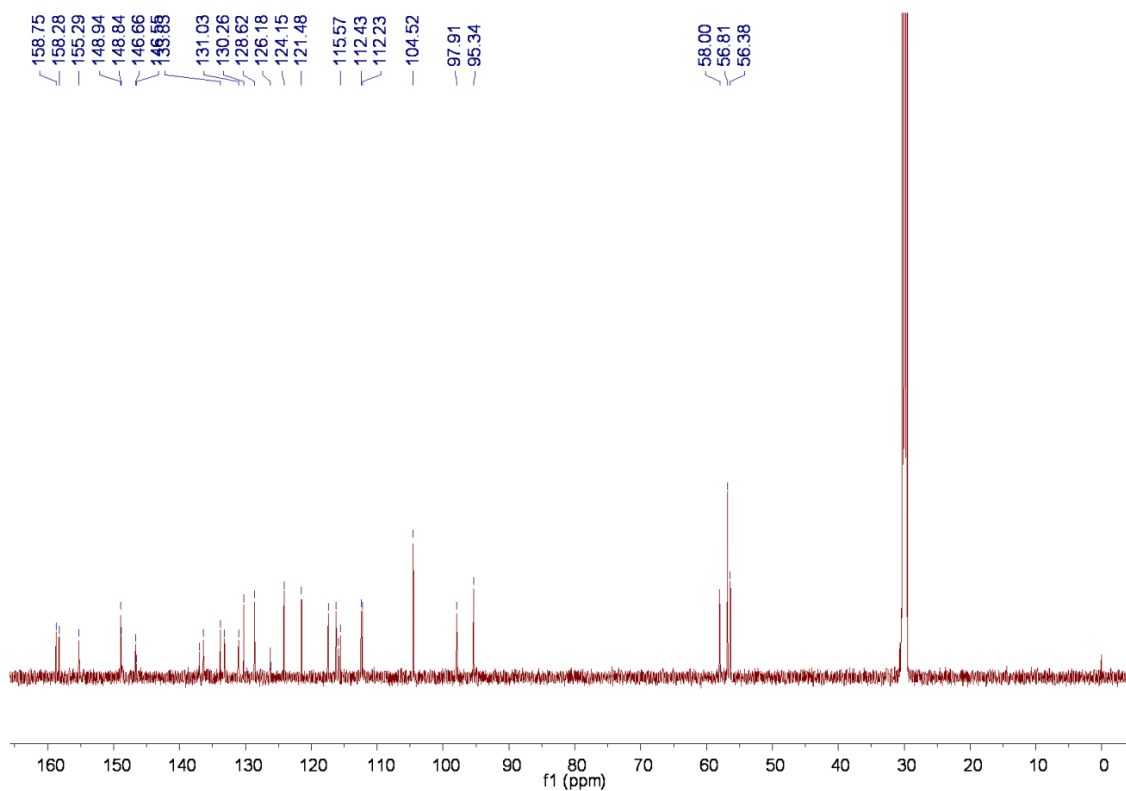
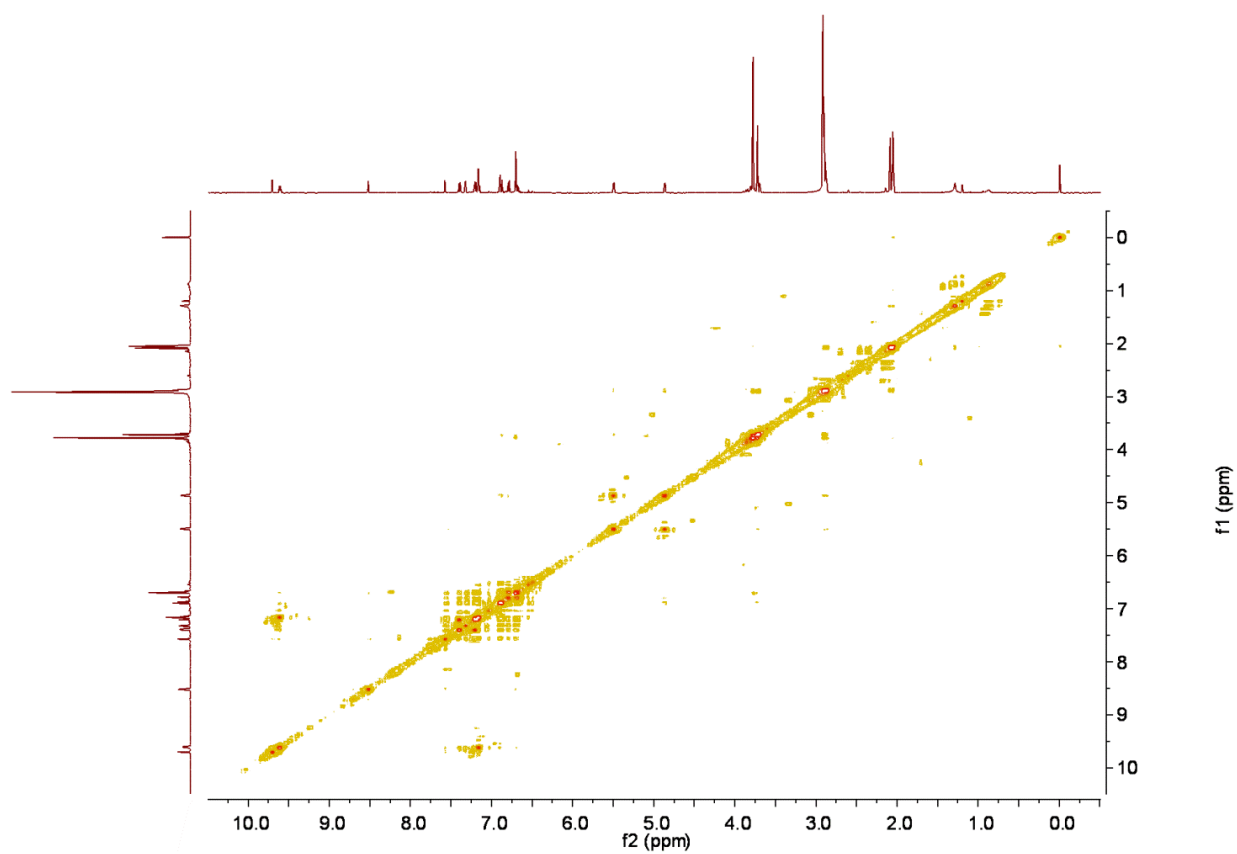
Figure S13. ^{13}C -NMR for loddigesinol I (3).**Figure S14.** ^1H - ^1H COSY for loddigesinol I (3).

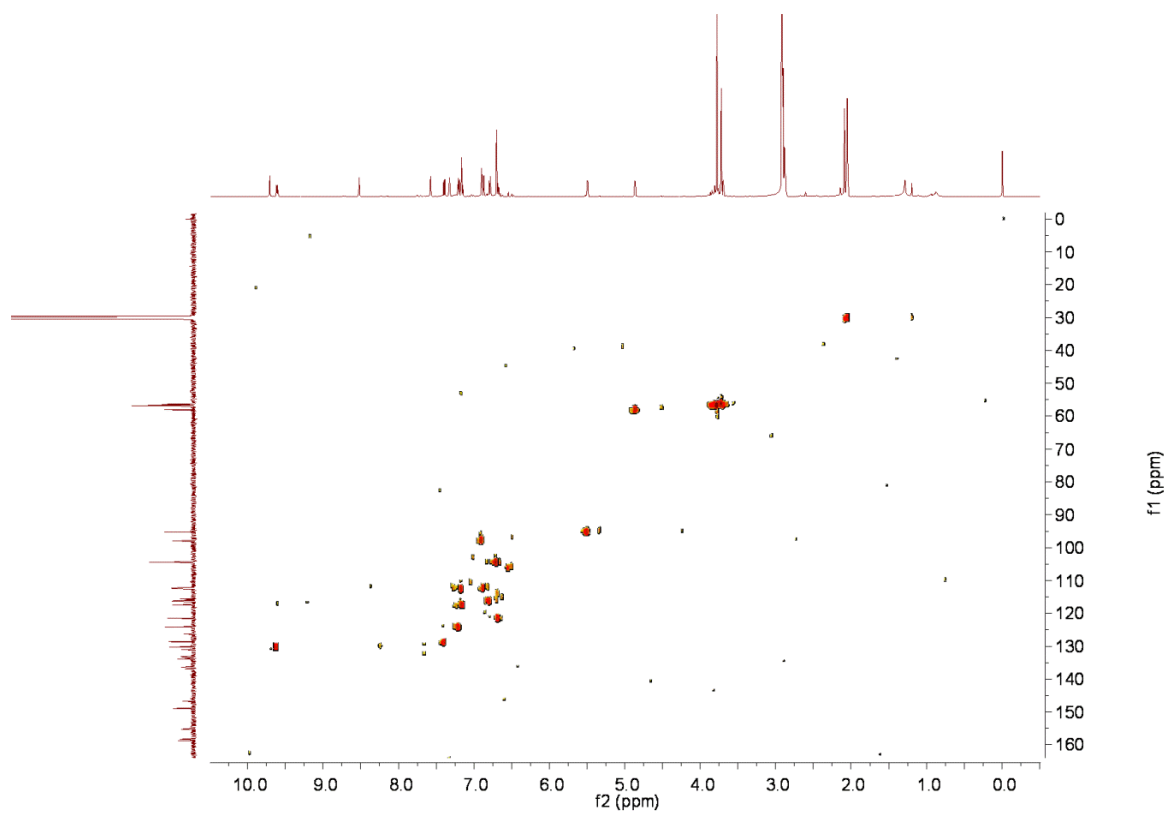
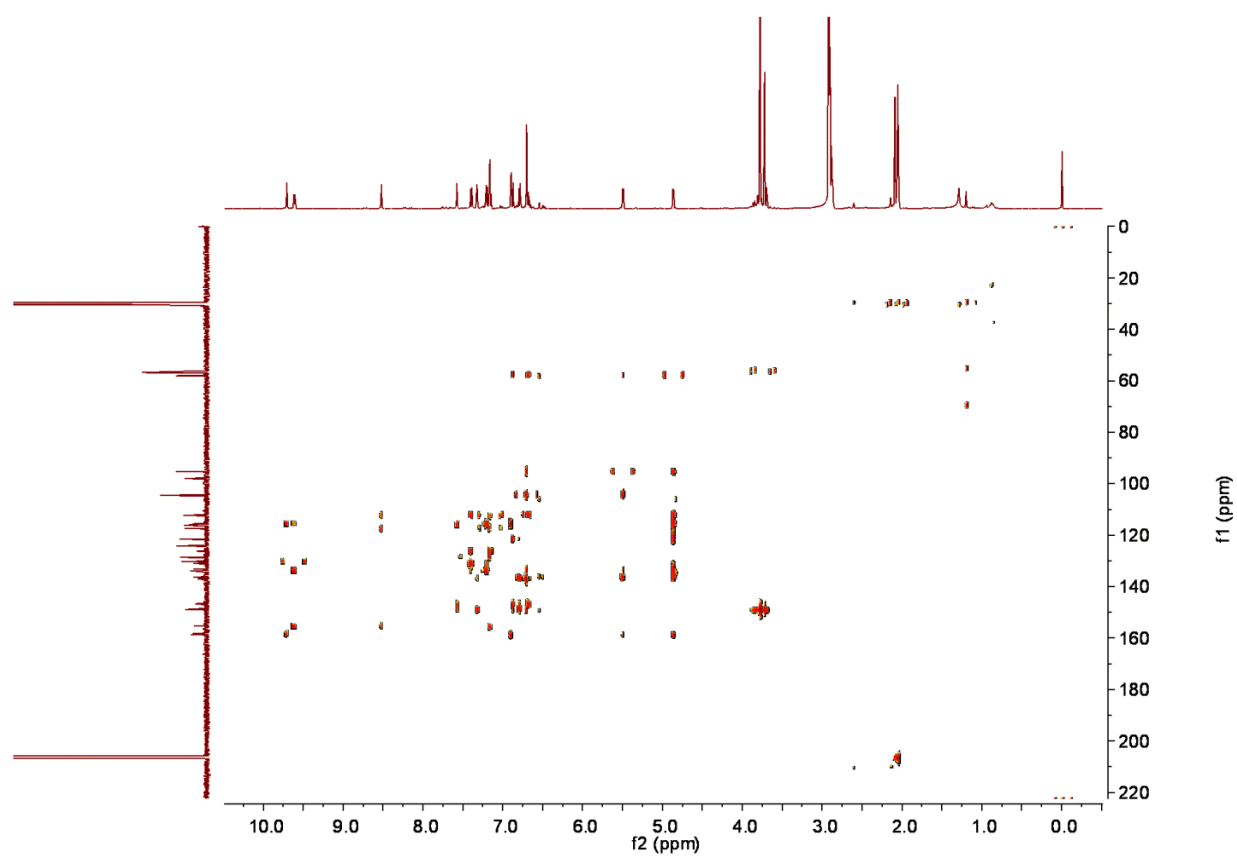
Figure S15. HSQC for loddigesiinol I (**3**).**Figure S16.** HMBC for loddigesiinol I (**3**).

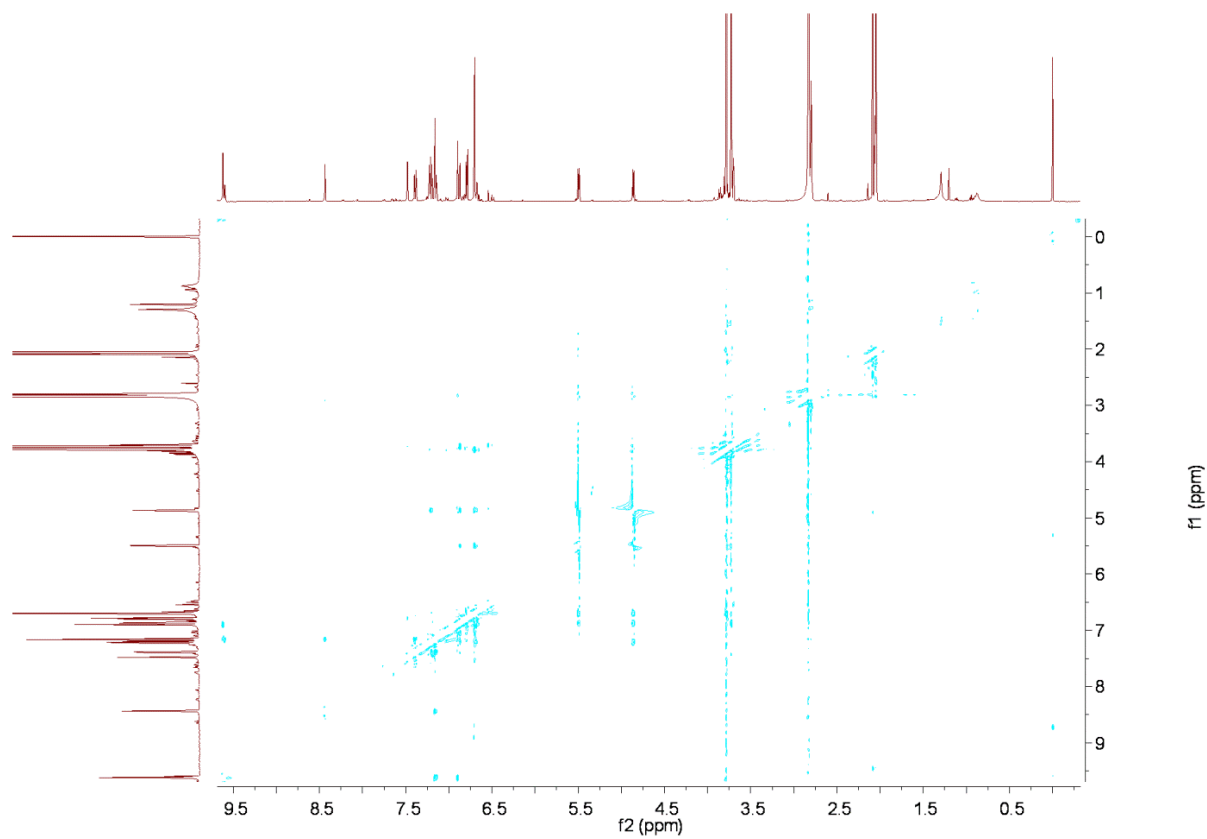
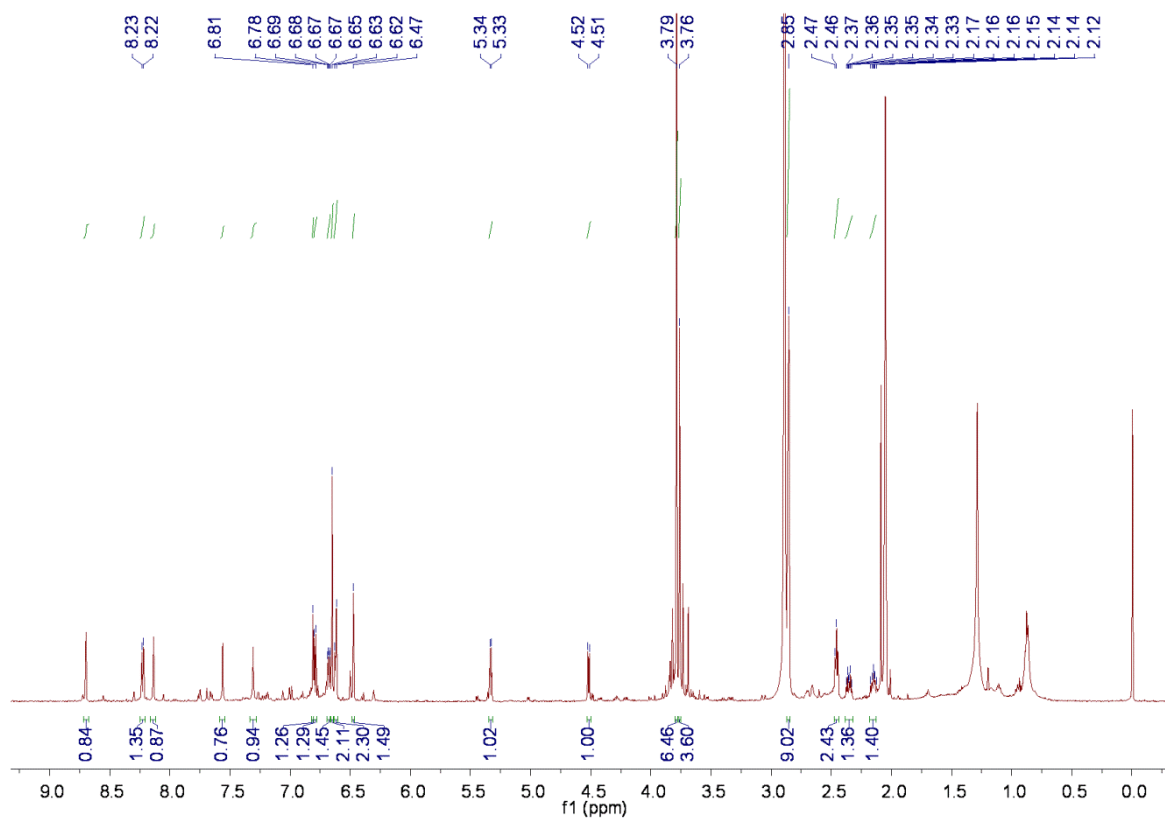
Figure S17. ROSEY for loddigesinol I (**3**).**Figure S18.** $^1\text{H-NMR}$ for loddigesinol J (**4**).

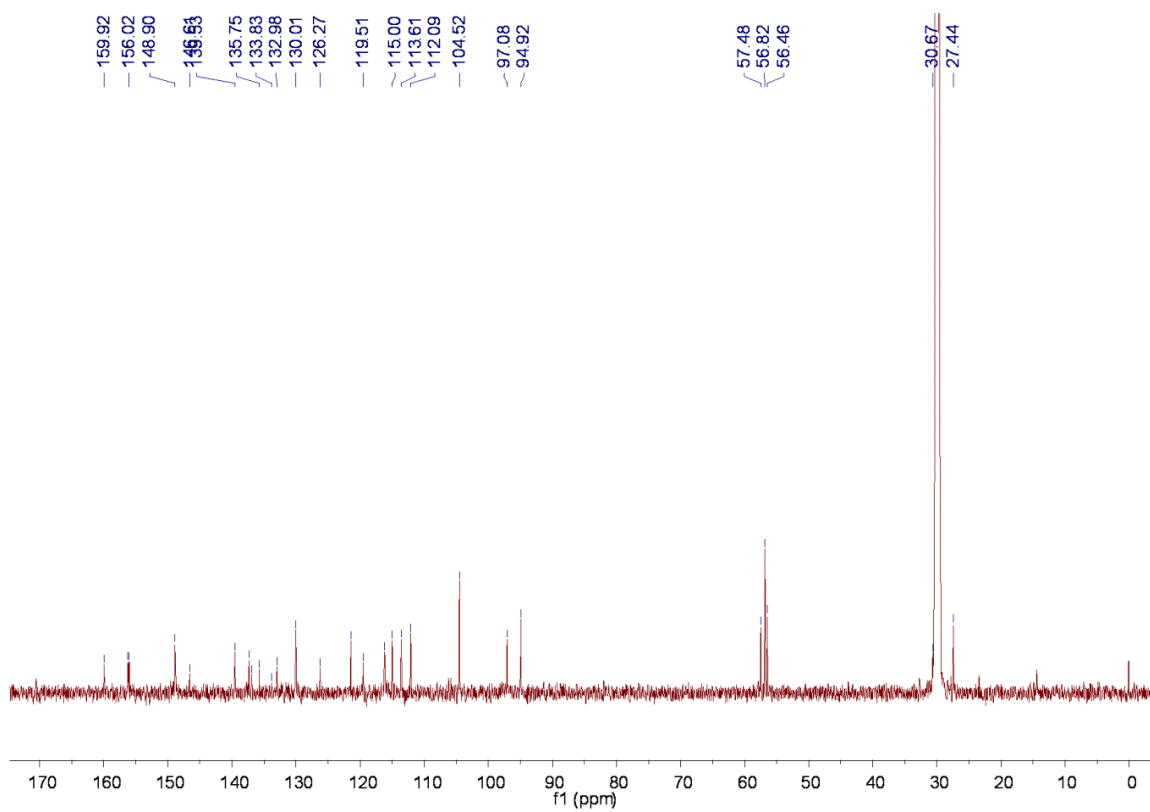
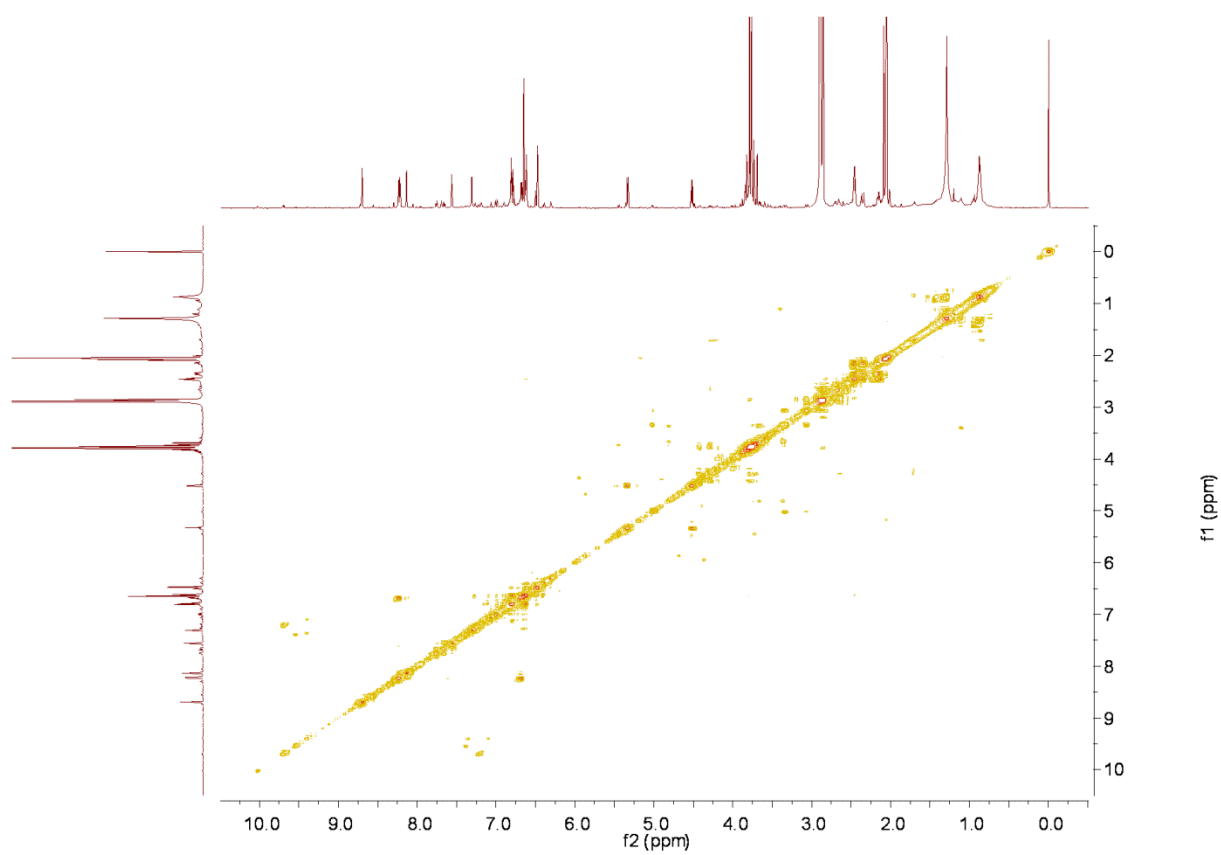
Figure S19. ^{13}C -NMR for loddigesinol J (4).**Figure S20.** ^1H - ^1H COSY for loddigesinol J (4).

Figure S21. HSQC for loddigesiinol J (4).

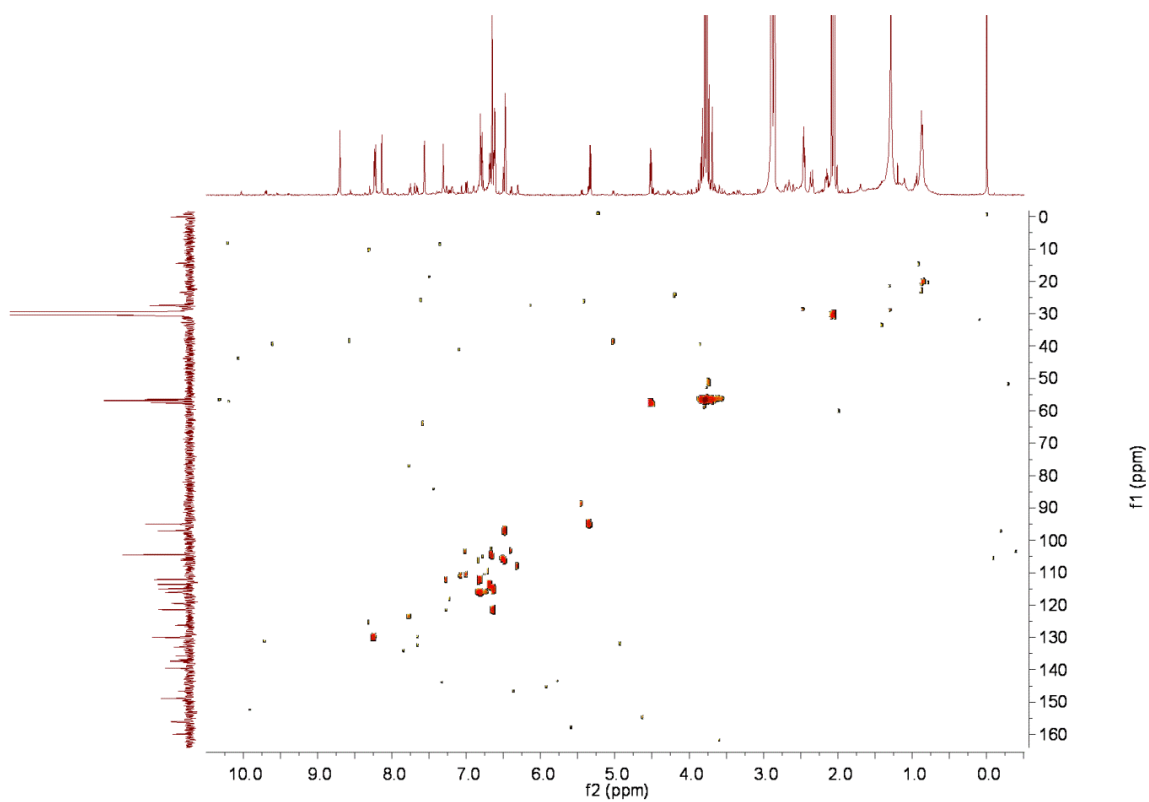


Figure S22. HMBC for loddigesiinol J (4).

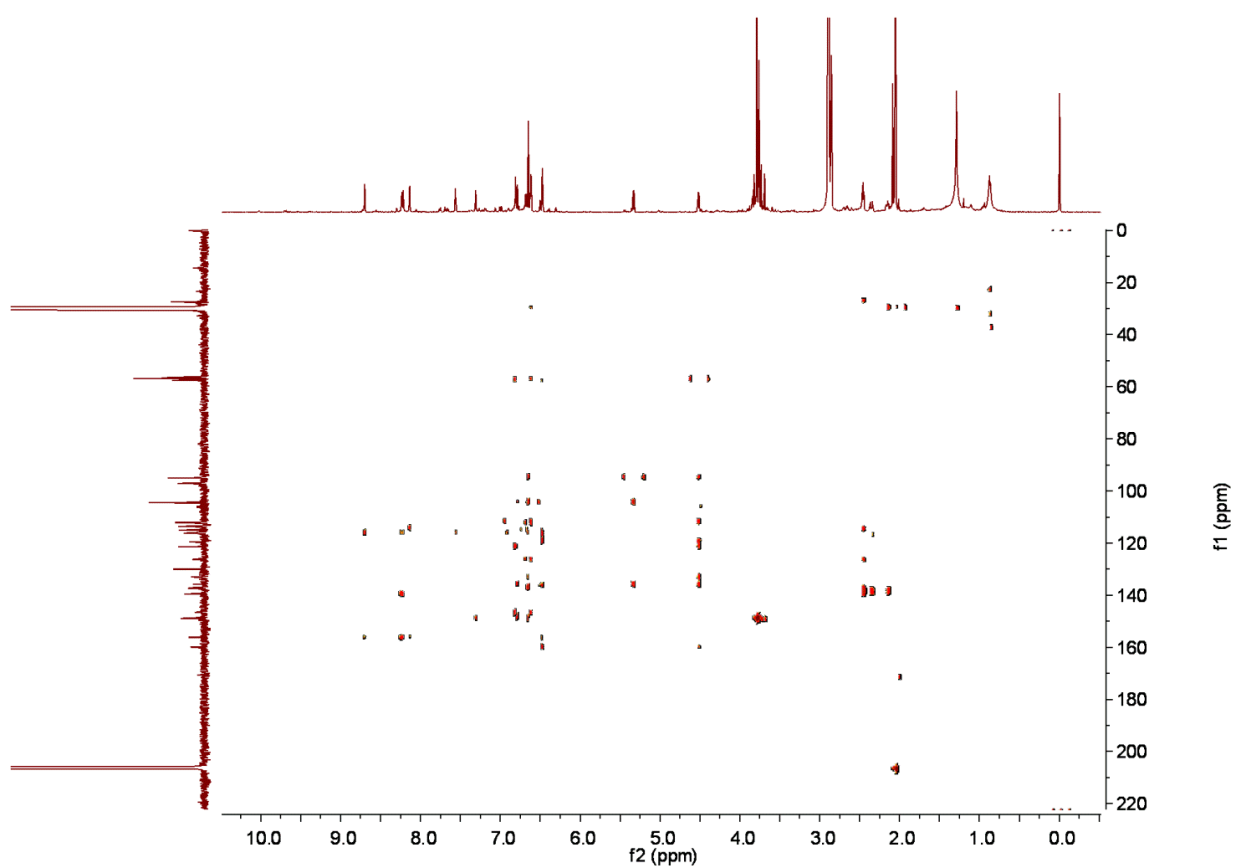


Figure S23. ECD spectrum for crepidatuol B (5).

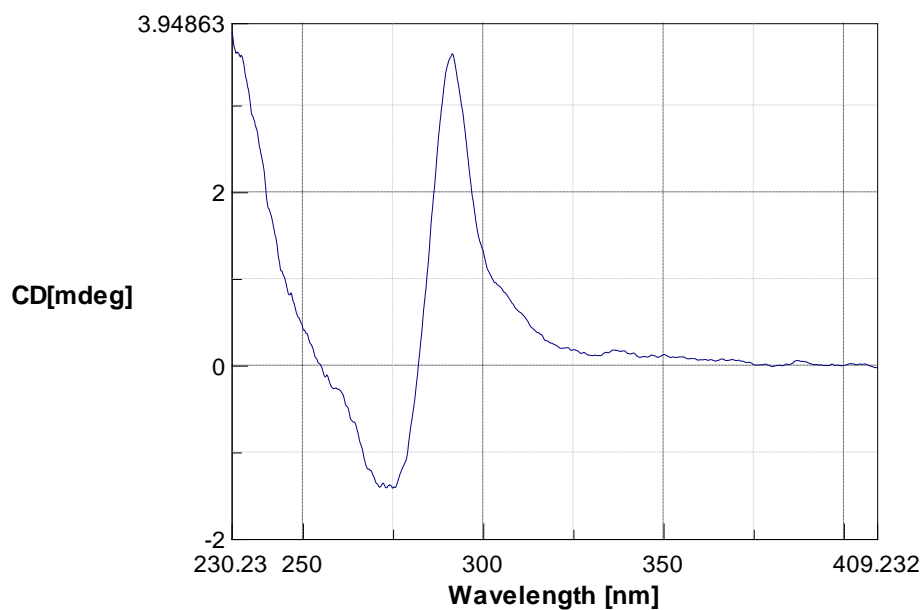


Figure S24. ^1H - ^1H COSY and HMBC correlations for compounds 1–5.

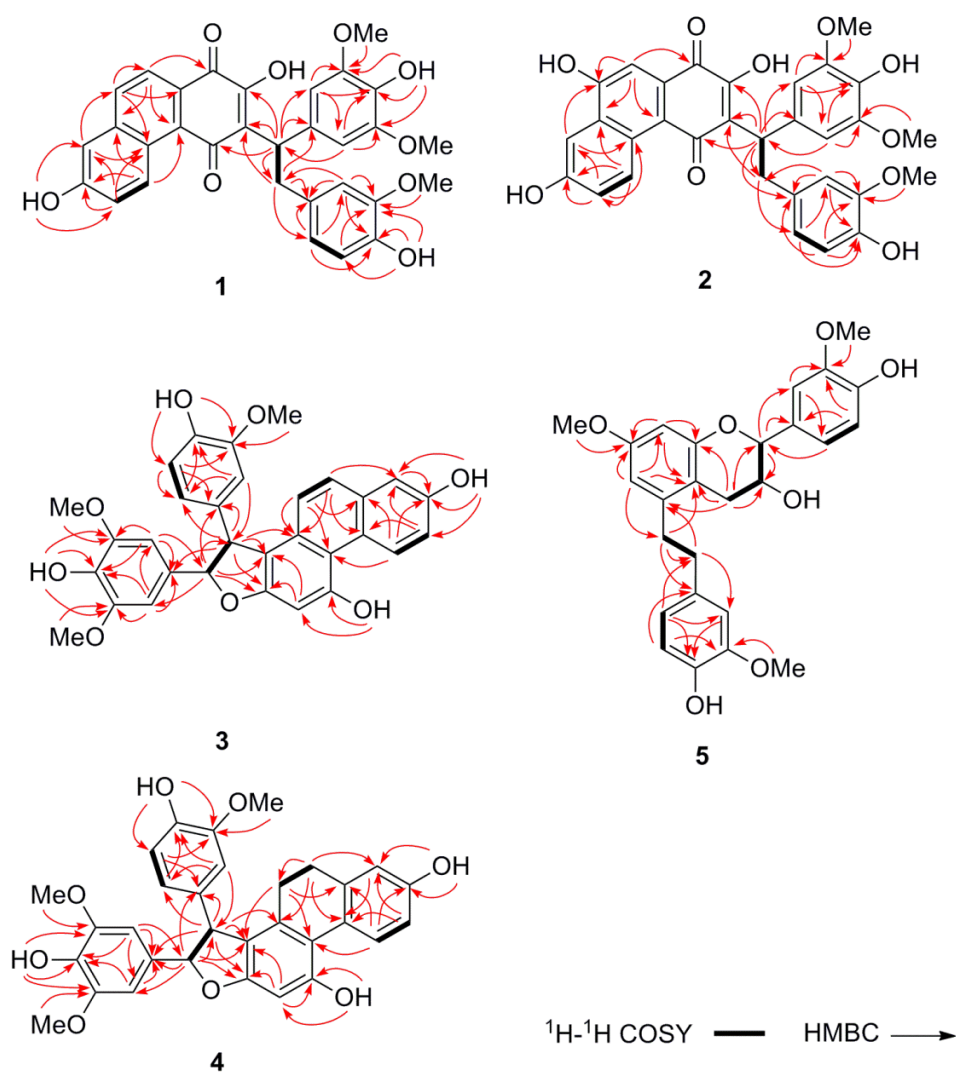


Table S1. ^{13}C Data for Compounds 1–4.

NO.	1^a	2^b	3^a	4^a
	δ	δ	δ	δ
1	183.0, C	183.4, C	115.6, C	119.5, C
2	153.4, C	152.9, C	158.7, C	159.9, C
3	126.8, C	126.4, C	97.9, CH	97.1, CH
4	189.5, C	188.6, C	158.3, C	156.2, C
4a	140.8, C	122.1, C	133.8, C	133.8, C
4b	125.4, C	127.7, C	126.2, C	126.3, C
5	131.6, CH	131.5, CH	130.3, CH	130.0, CH
6	123, CH	122.8, CH	117.4, CH	113.6, CH
7	158.7, C	157.9, C	155.3, C	156.0, C
8	111, CH	105.4, CH	112.4, CH	115.0, CH
8a	129.7, C	131.4, C	115.8, C	139.5, C
9	132.9, CH	157.2, C	128.6, CH	30.2, CH ₂
10	122.3, CH	103.8, CH	124.1, CH	27.5, CH ₂
10a	128.1, C	129.8, C	131.0, C	136.9, C
1''	134.6, C	134.8, C	133.2, C	133.0, C
2''	107.4, CH	107.4, CH	104.5, CH	104.5, CH
3''	148.5, C	148.4, C	148.9, C	148.9, C
4''	135.6, C	135.5, C	136.9, C	137.3, C
5''	overlap with 3'	overlap with 3'	overlap with 3'	overlap with 3'
6''	overlap with 2'	overlap with 2'	overlap with 2'	overlap with 2'
a	133.2, C	133.3, C	136.4, C	135.8, C
a'	113.4, CH	113.3, CH	112.2, CH	112.1, CH
3 ¹ -CH ₃ O	148.0, C	147.9, C	148.8, C	148.7, C
3 ^{''} -CH ₃ O	145.7, C	145.6, C	146.6, C	146.6, C
5 ¹ -CH ₃ O	115.6, CH	115.5, CH	116.2, CH	116.2, CH

The data were recorded at ^a150 MHz and ^b125 MHz; chemical shifts (δ) are in ppm.

Table S2. ¹H-NMR and HMBC Data for Compounds 1–4.

1 ^a		2 ^b		3 ^a		4 ^a		
No.	δ(J)	HMBC	δ(J)	HMBC	δ(J)	HMBC	δ(J)	HMBC
3					6.9, s	C-1, 2	6.47, s	C-1, 2, 4
5	9.39, d (9.6)	C-4a, 7, 8, 8a	9.45, d (9.5)	C-6, 7, 8, 8a, 4a	9.61, d (9.0)	C-8a, 4a, 7	8.22, d (9.0)	C-7, 8a
6	7.35, d (9.6)	C-4b, 7, 8	7.33, dd (9.5, 3.0)	C-4b, 7, 8	7.15, dd (9.0, 2.4)	C-4b, 7, 8	6.67, dd (8.4, 1.8)	C-4b, 8
8	7.26, brs	C-4b, 7, 9	7.63, brs	C-4b, 6, 9	7.16, brs	C-9, 7, 6	6.62, brs	C-4b, 9, 6
9	7.96, d (8.4)	C-4a, 4b, 8, 10, 10a			7.39, d (9.0)	C-4b, 8, 10a, 4a	2.46, t (7.8)	C-4b, 8, 10, 8a
10 _α	7.93, d (8.4)	C-1, 4, 4b, 4a, 8a	7.38, s	C-1, 4, 4a, 8, 8a, 9, 10a	7.20, d (9.6)	C-8a, 10a	2.35, td (15.6, 7.8)	C-1, 9, 8a
10 _β							2.15, dt (15.0, 7.8)	C-1, 9, 8a
2'	6.90, s	C-3', 4', 5', 6', a	6.89, s	C-3', 4', 5', 6', a	6.70, s	C-1', 6', 3', 4', a	6.65, s	C-1', 6', 3', 4', a
6'	6.90, s	C-2', 3', 4', 5', a	6.89, s	C-2', 3', 4', 5', a	6.70, s	C-1', 2', 3', 4', a	6.65, s	C-1', 2', 5', 4', a
2''	6.85, s	C-1'', 3'', 4'', 6'', a'	6.85, d (2.0)	C-1'', 3'', 4'', 5'', 6'', a'	6.87, s	C-4'', 6'', a'	6.81, s	C-4'', 6'', a'
5''	6.60, d (7.8)	C-1'', 3''	6.61, d (8.0)	C-1'', 2'', 3'', 4'', 6''	6.79, d (8.4)	C-1'', 3'', 6''	6.79, d (7.8)	C-1'', 3''
6''	6.71, d (7.8)	C-2'', 4'', a'	6.70, dd (8.0, 2.0)	C-2'', 4'', a'	6.67, d (7.8)	C-2'', 4'', a'	6.62, d (7.8)	C-2'', 4'', a'
a	4.81, dd (10.2, 6.6)	C-2, 3, 4, 1', 2'/6', a'	4.80, dd (10.0, 6.5)	C-2, 3, 4, 1', 2', a'	5.49, d (6.6)	C-2, 2'/6', 1'', a', 1	5.33, d (7.2)	C-2, 1', 2'/6', 1'', a'
a' _α	3.36, dd (13.8, 6.6)	C-1', 1'', 2'', 6'', a	3.67, dd (13.5, 10.0)	C-3, 1'', 2'', 6'', a	4.86, d (6.6)	C-2, 1', 1'', 6'', a, 10a, 1	4.52, d (6.6)	C-1, 2, 1', 1'', 2''/6'', a, 10a
a' _β	3.66, dd (13.8, 10.2)	C-1', 1'', 2'', 6'', a	3.34, dd (13.5, 7.0)	C-3, 1'', 2'', 6'', a				
3'-CH ₃ O	3.80, s	C-3'	3.8, s	C-3'	3.78, s	C-3', 5'	3.79, s	C-3', 5'
3''-CH ₃ O	3.67, s	C-3''	3.67, s	C-3''	3.72, s	C-3''	3.76, s	C-3''

Table S2. Cont.

No.	1^a		2^b		3^a		4^a	
	δ (J)	HMBC	δ (J)	HMBC	δ (J)	HMBC	δ (J)	HMBC
5'-CH ₃ O	overlap with 3'	C-5'	overlap with 3'	C-5'	overlap with 3'	overlap with 3'	overlap with 3'	overlap with 3'
9-OH			unobserved					
2-OH	unobserved		unobserved					
7-OH	9.36, brs	C-6, 8	unobserved		8.52, s	C-7, 6, 8	8.70, s	C-7, 8
4-OH					9.7, s	C-4, 3	8.13, s	C-4, 3
4''-OH	7.21, brs		unobserved		7.57, s	C-3'', 5''	7.56, s	C-3'', 5''
4'-OH	7.02, brs		unobserved		7.32, s	C-3', 5', 4'	7.31, s	C-3', 5', 4'

The data were recorded at ^a 600 MHz and ^b 500 MHz; chemical shifts (δ) are in ppm, and coupling constants (J) are in Hz.

Table S3. NMR Data of Compounds **5** (in C₂D₆CO).

5^d			
NO.	δ_H (J)	HMBC	δ_c
1			82.9, CH
2	4.60, d (8.0)	C-3, 1', 2', 6', 4	68.6, CH
3	4.05, m		31.8, CH ₂
4 _{α}	3.00, dd (15.6, 5.2)	C-3, 2, 5, 9, 10	
4 _{β}	2.66, dd (15.6, 8.8)	C-3, 2, 5, 9, 10	143.1, C
5			109.2, CH
6	6.42, d (2.4)	C-7, 8, a', 10	159.9, C
7			99.9, CH
8	6.27, d (2.4)	C-9, 7, 6	156.5, C
9			112.5, C
10			131.8, C
1'			111.9, CH
2'	7.04, s	C-2, 3', 6'	147.5, C
3'			143.1, C
4'			115.6, CH
5'	6.82, d (8.0)	C-1', 3'	121.5, CH
6'	6.89, d (8.0)	C-2, 2', 3'	134.1, C
1''			113.1, CH
2''	6.83, s	C4'', 6'', a	148.2, C
3''			145.7, C
4''			115.7, CH
5''	6.74, d (8.0)	C-1'', 3''	121.7, CH
6''	6.69, d (8.0)	C-2'', a, 4''	37, CH ₂
a	2.81, m	C-5, 1'', 6'', 10, a'	36.1, CH ₂
a'	2.81, m	C-5, 1'', 6'', 10, a'	56.4, CH ₃
3'-CH ₃ O	3.85, s	C-3'	56.3, CH ₃
3''-CH ₃ O	3.81, s	C-3''	55.5, CH ₃
7-CH ₃ O	3.72, s	C-7	
3-OH	4.00, brs		
4'-OH	7.60, brs		
4''	7.31, brs		

The data were recorded at ^d 400 MHz (¹H-NMR) and 100 MHz (¹³C-NMR); chemical shifts (δ) are in ppm and coupling constants (J) are in Hz.