

Supporting Information

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Myrrhalindenane C, A New Eudesmane Sesquiterpenoid From *Lindera Myrrha* Roots

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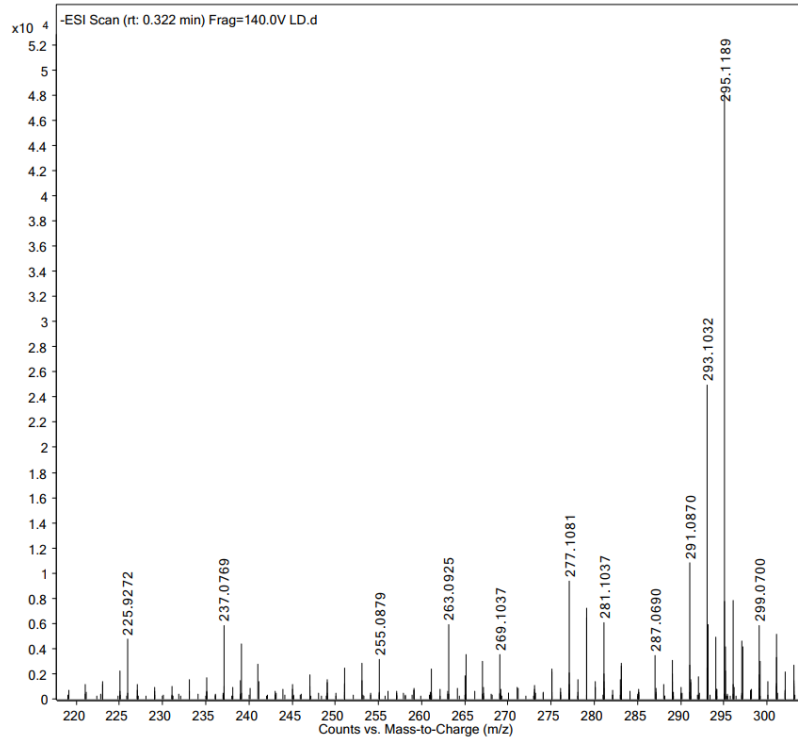


Figure S1: HRESIMS spectrum of 1.

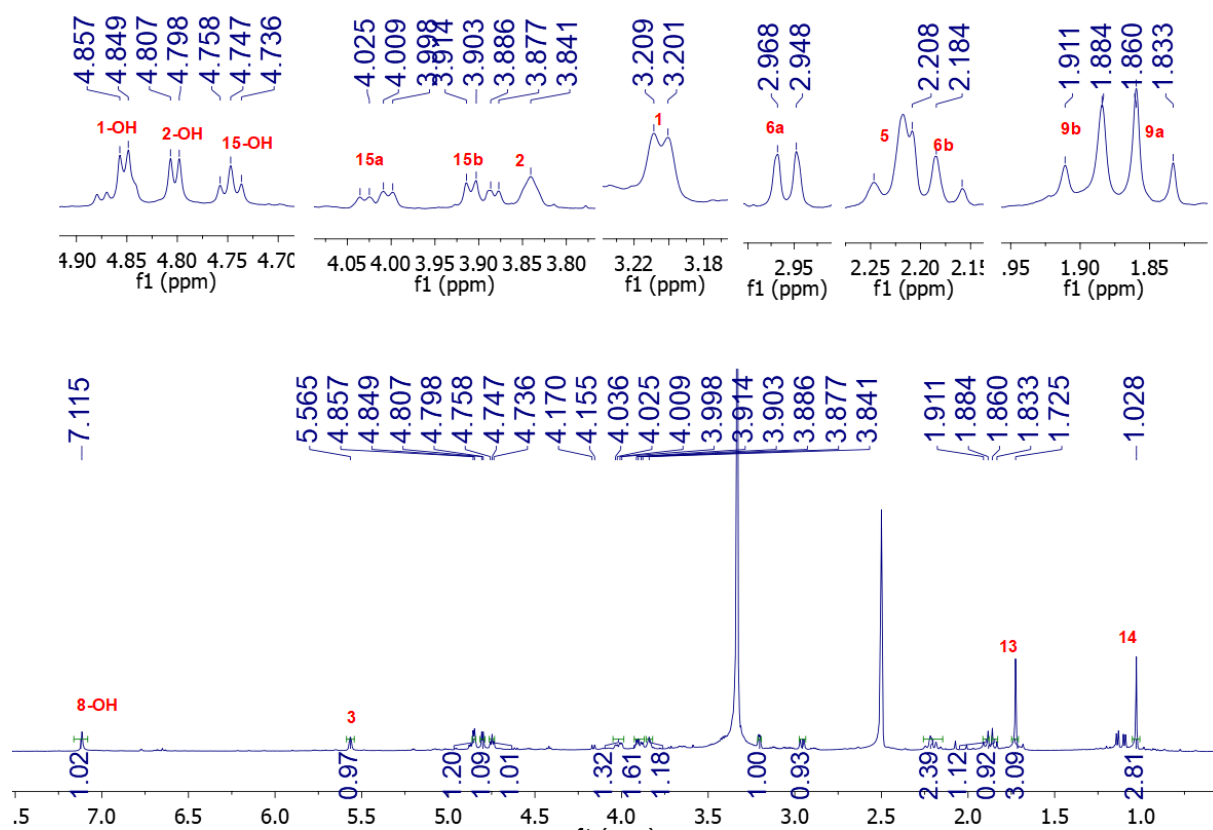


Figure S2: The ^1H NMR spectrum of **1** in $\text{DMSO-}d_6$.

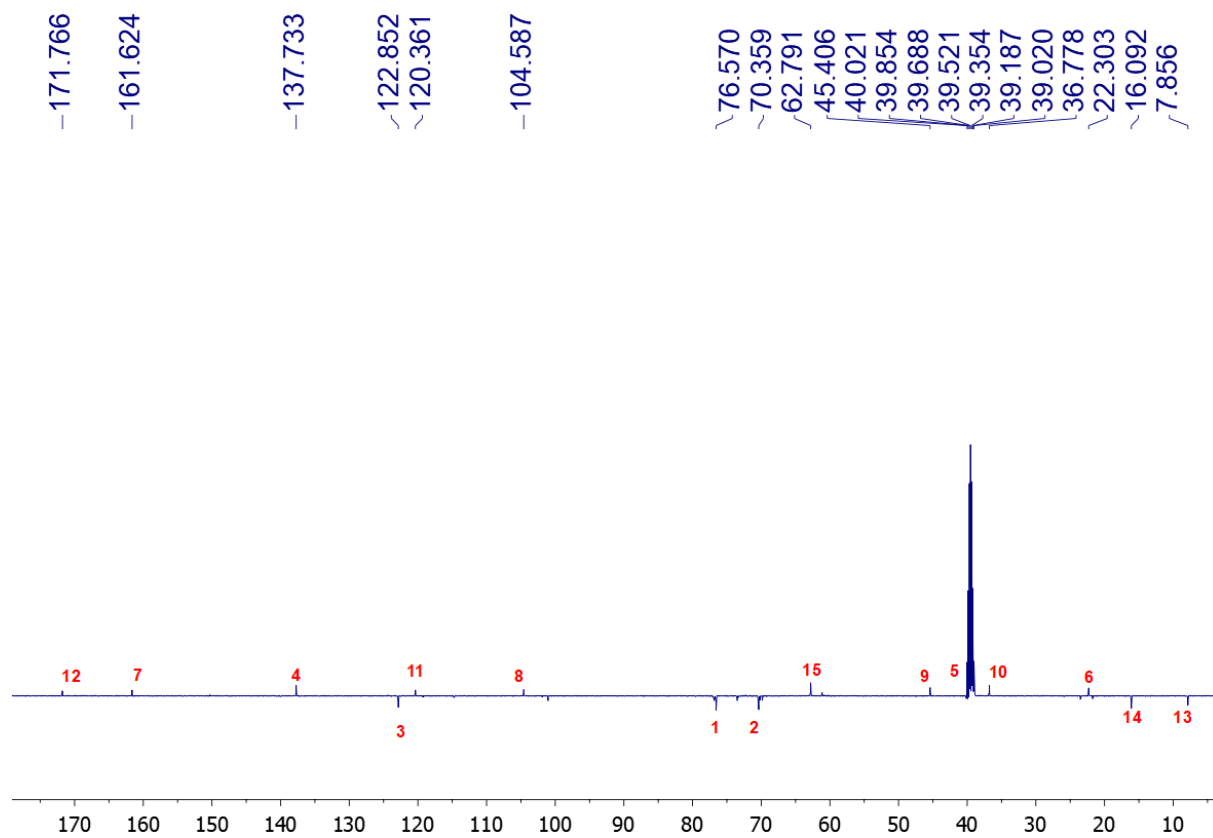


Figure S3: The ^{13}C NMR spectrum of **1** in $\text{DMSO-}d_6$.



Figure S4: The HSQC spectrum of **1** in DMSO- d_6 .

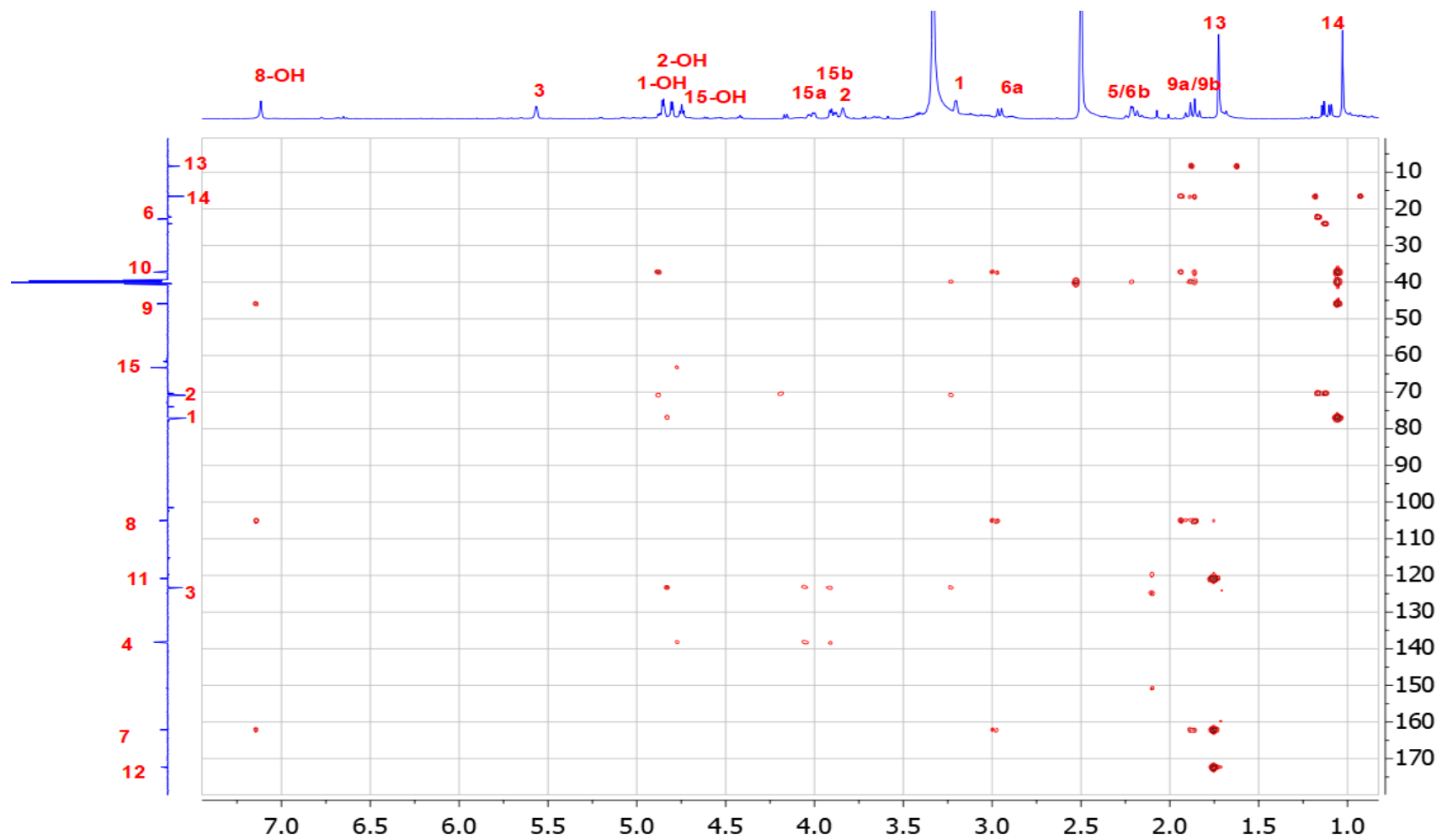


Figure S5: The HMBC spectrum of **1** in DMSO-*d*₆.

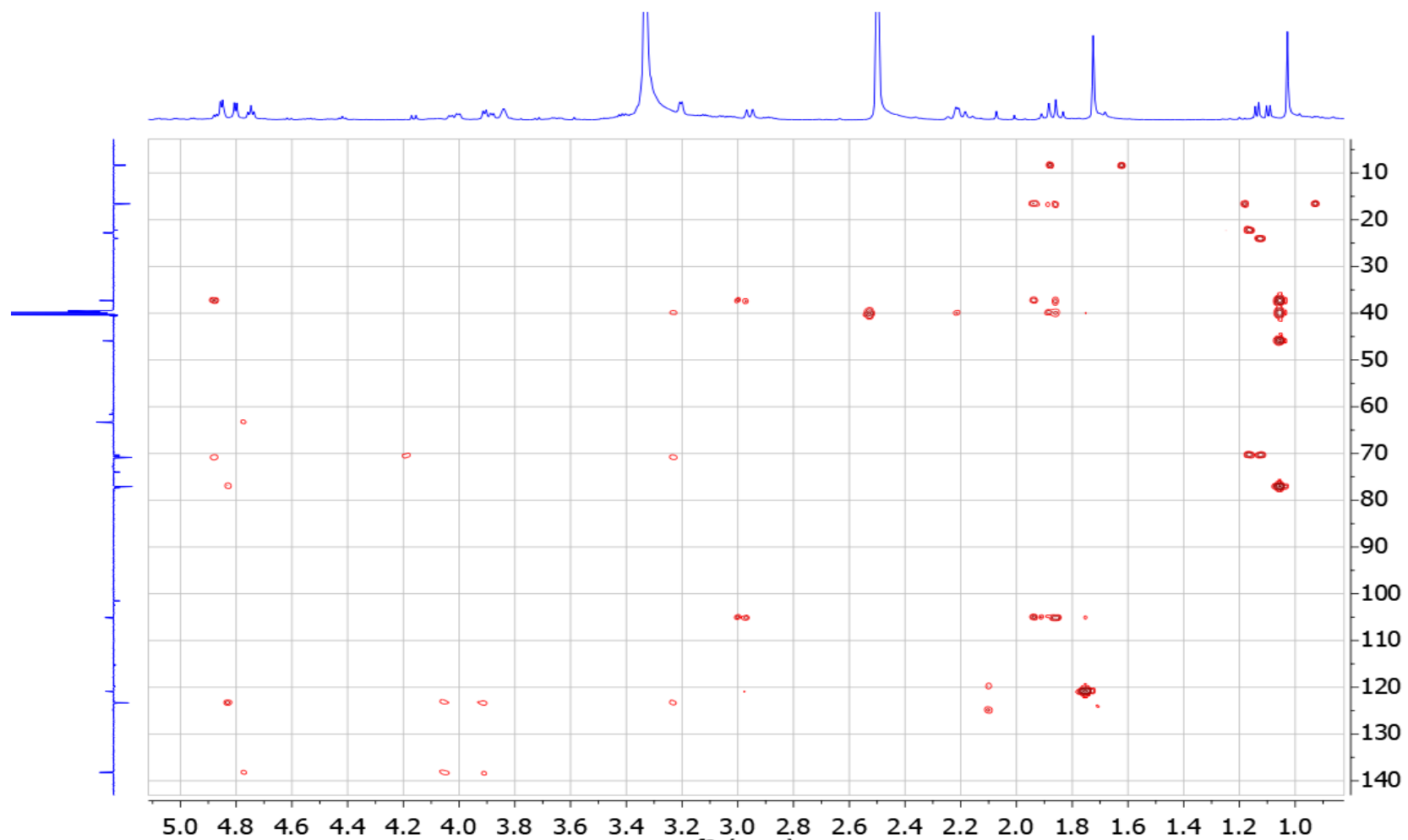


Figure S6: The HMBC spectrum of **1** in DMSO-*d*₆ (expansion).

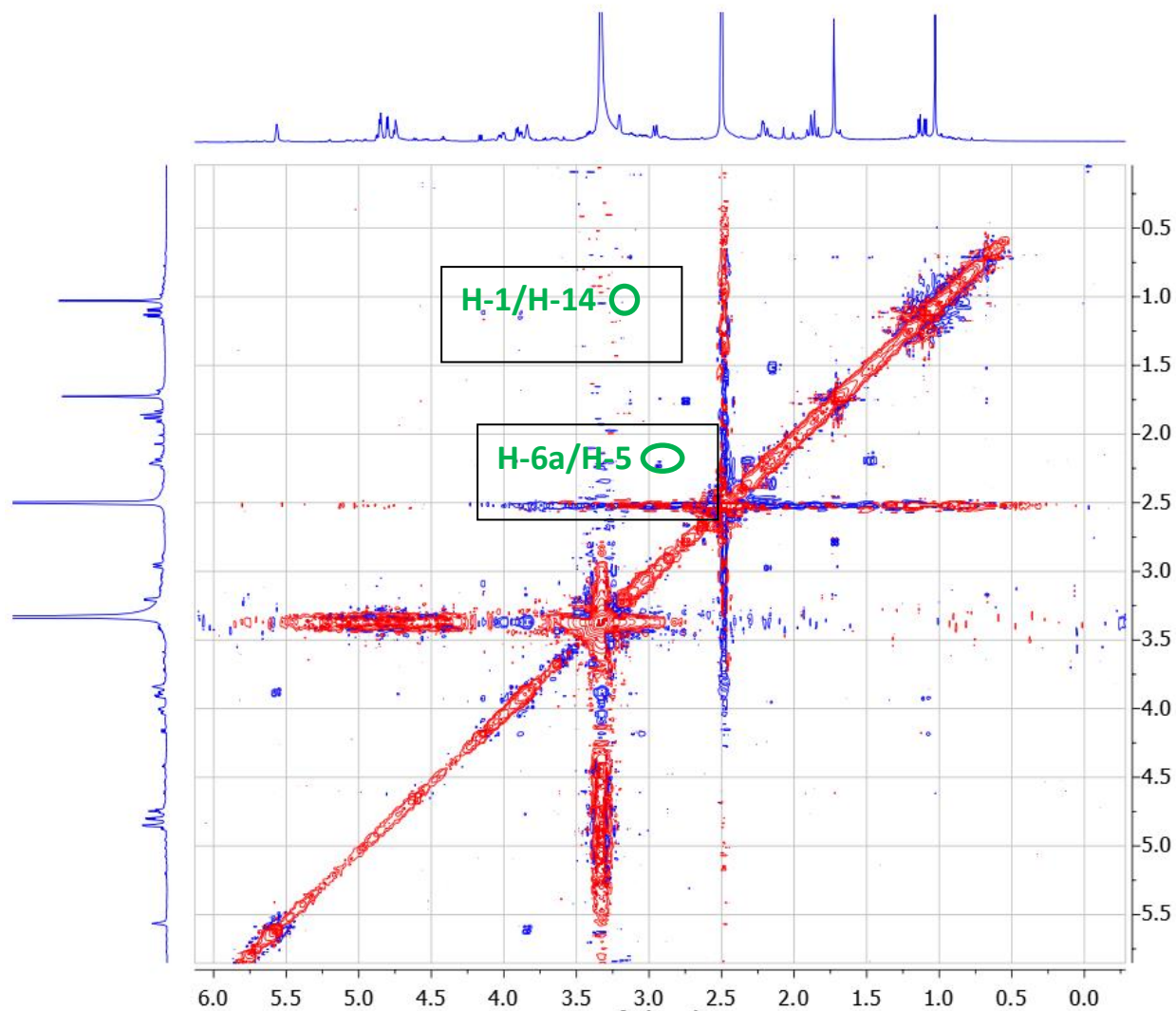


Figure S7: The NOESY spectrum of **1** in DMSO-*d*₆.

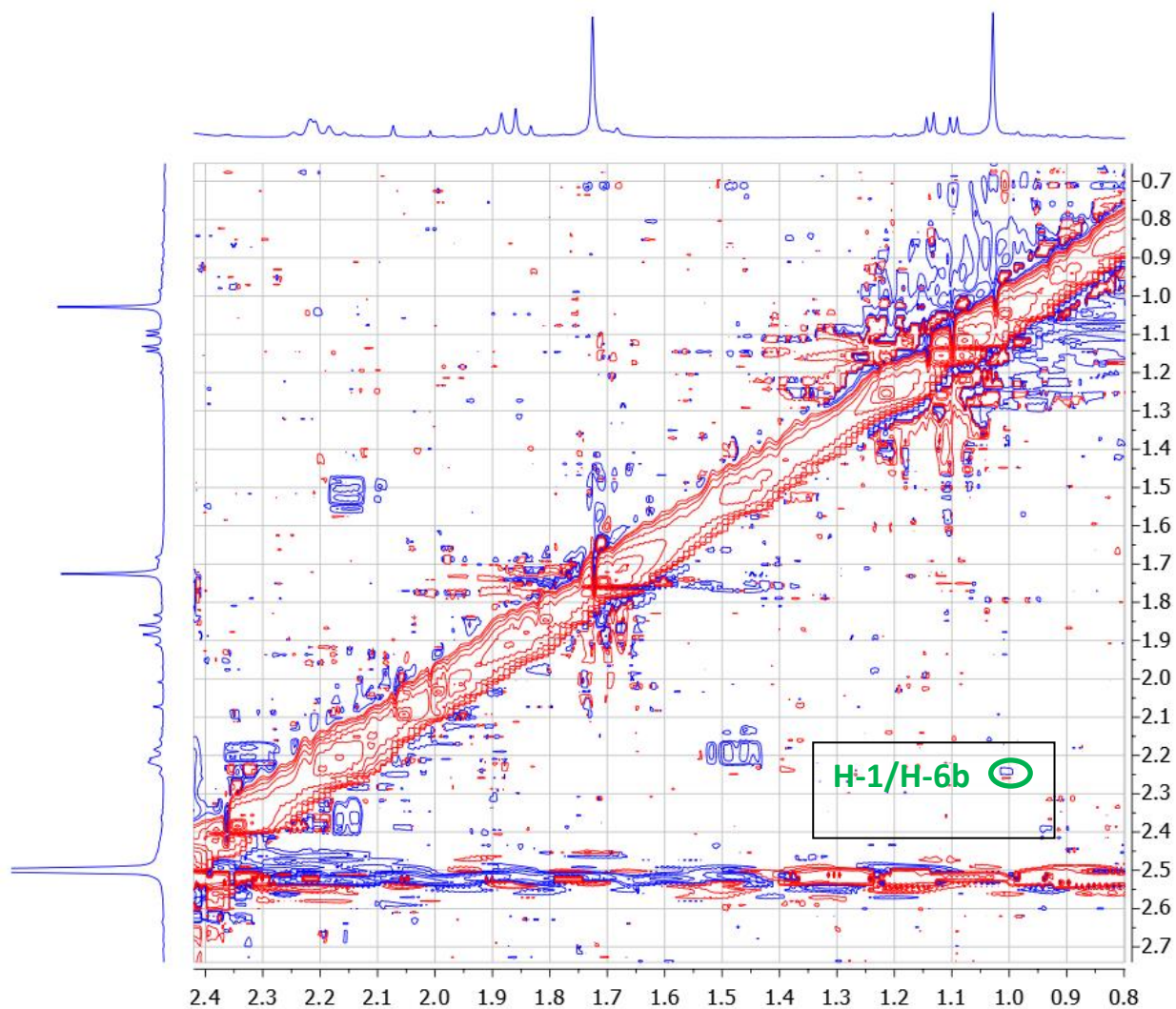


Figure S8: The NOESY spectrum of **1** in DMSO-*d*₆ (expansion).

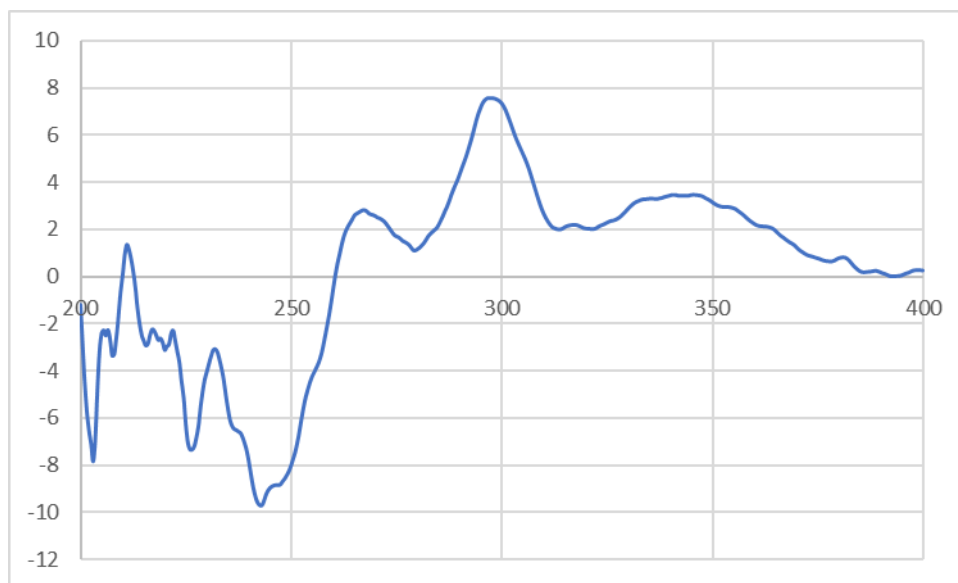


Figure S9: ECD spectrum of **1**.

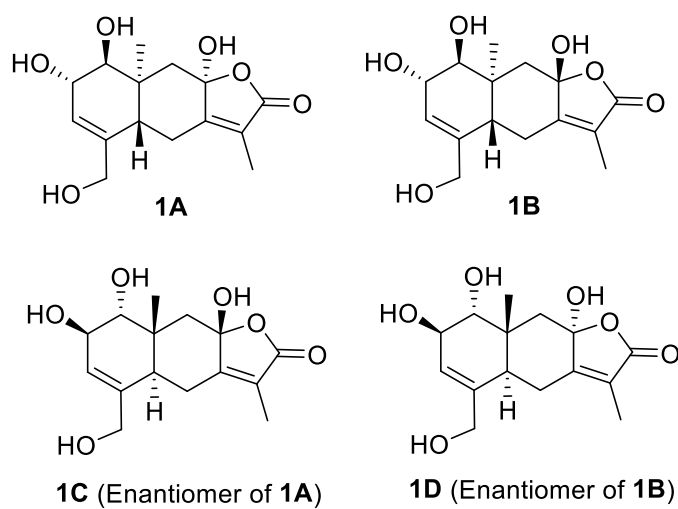


Figure S10: Four possible stereoisomers of **1**.

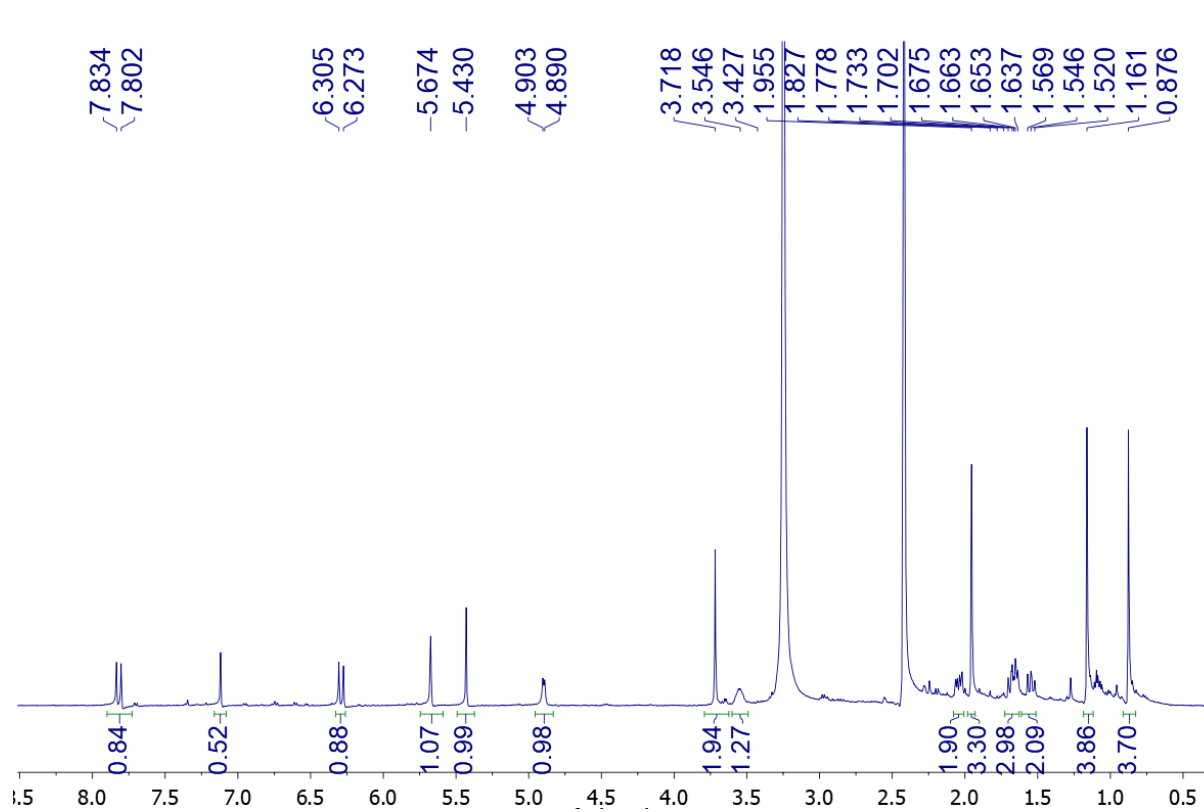


Figure S11: The ^1H NMR spectrum of **2** in $\text{DMSO-}d_6$.

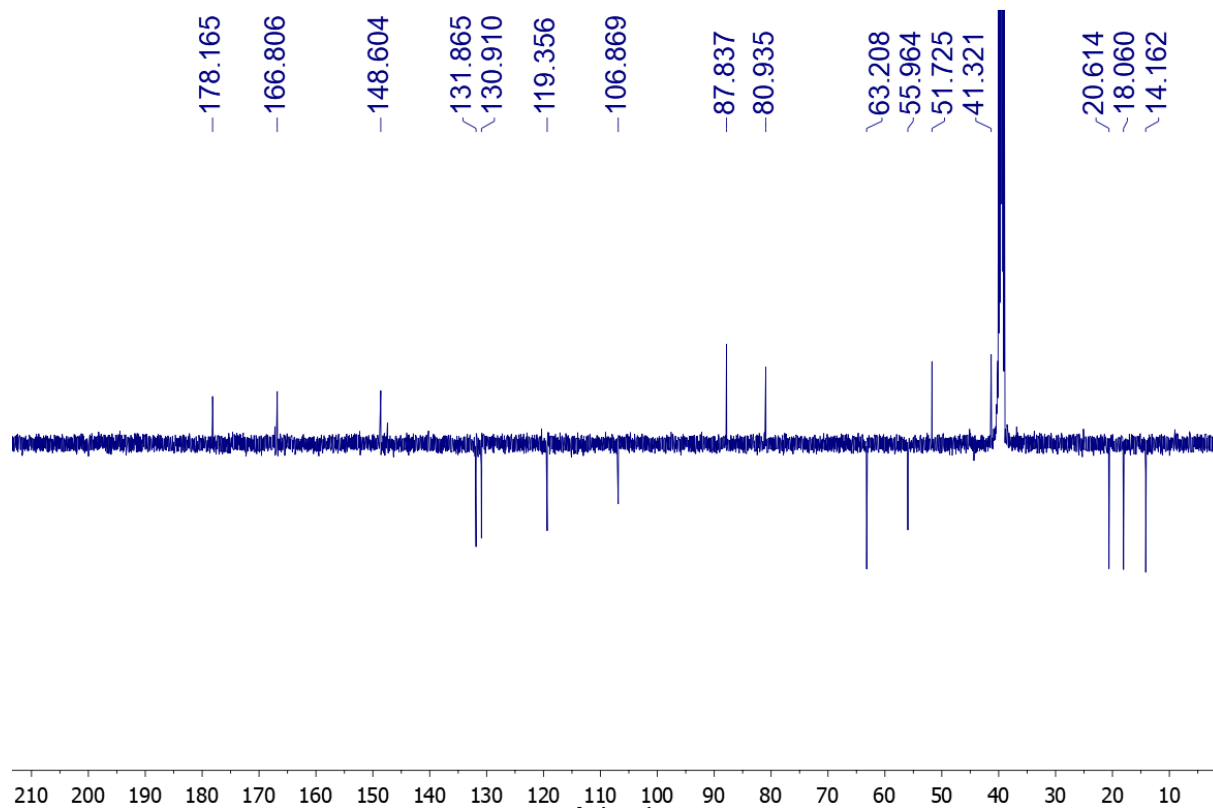


Figure S12: The ^{13}C NMR spectrum of **2** in $\text{DMSO-}d_6$.



Figure S13: HMBC spectrum of **2** in DMSO-*d*₆.

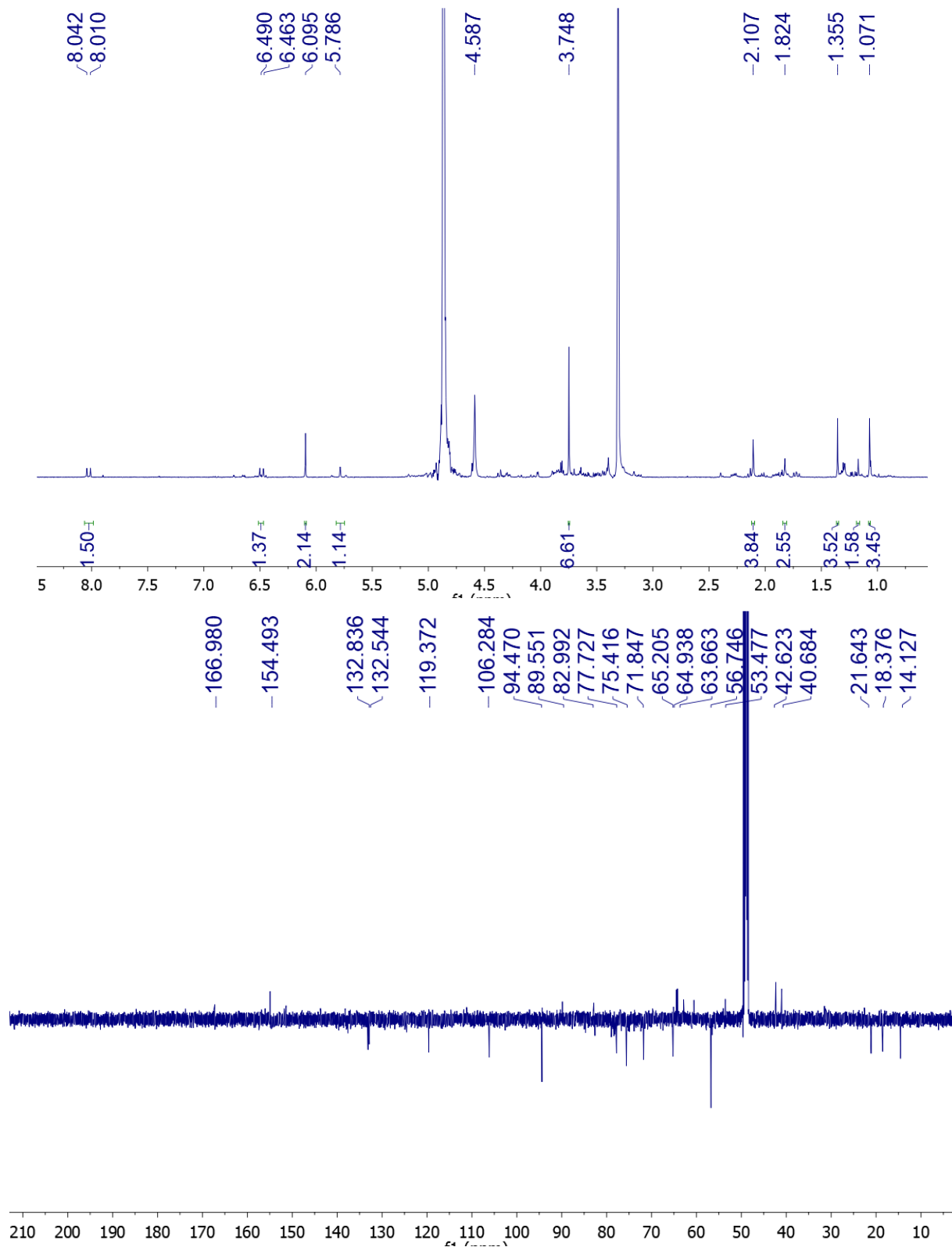


Figure S14: The ¹H and ¹³C NMR spectrum of **3** in methanol-*d*₄

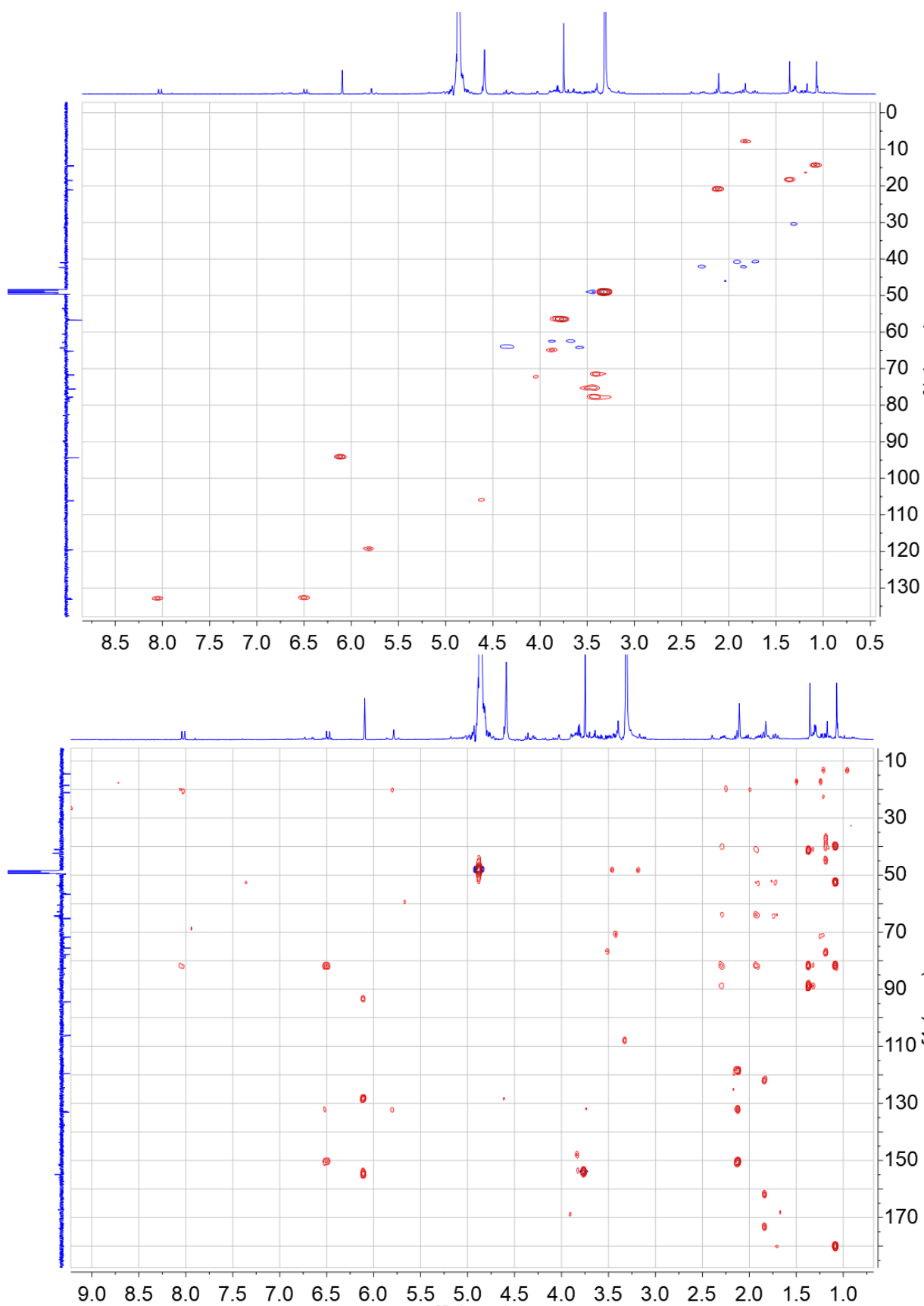


Figure S15: HSQC and HMBC spectra of **3** in methanol-*d*₄

OD2.30

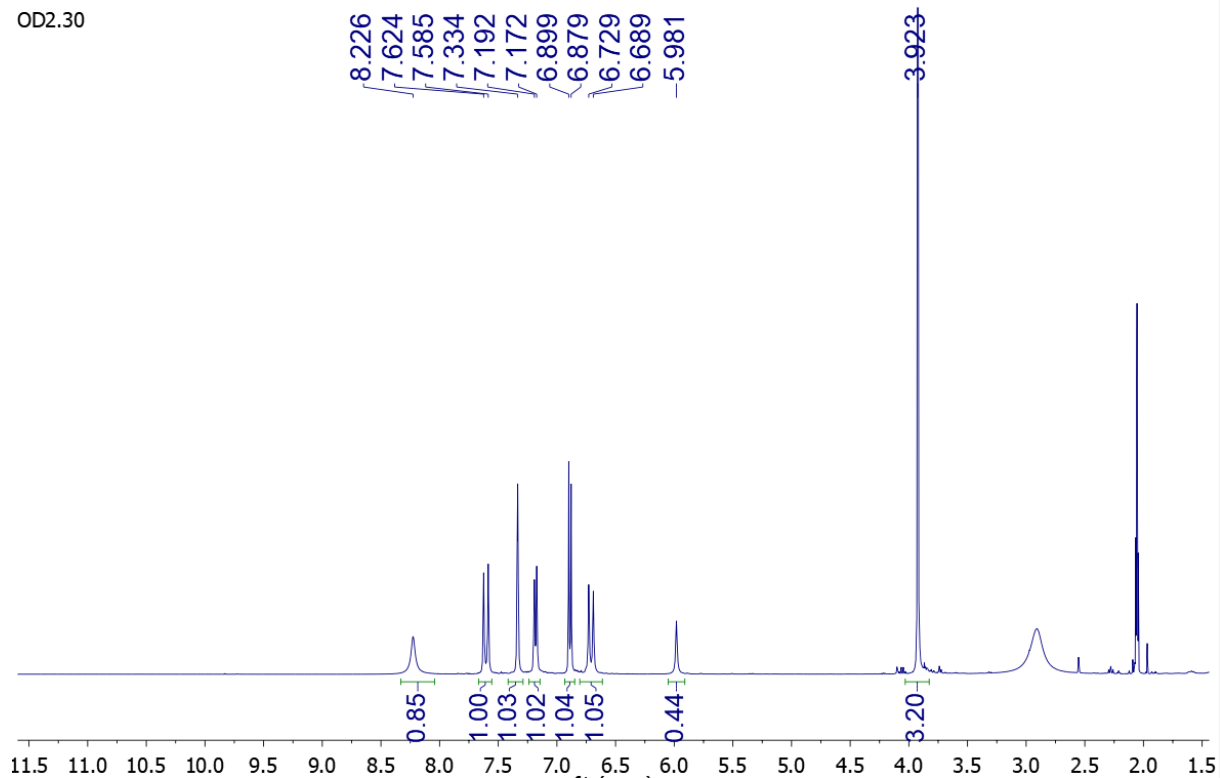


Figure S16: The ^1H NMR spectrum of **4** in acetone- d_6

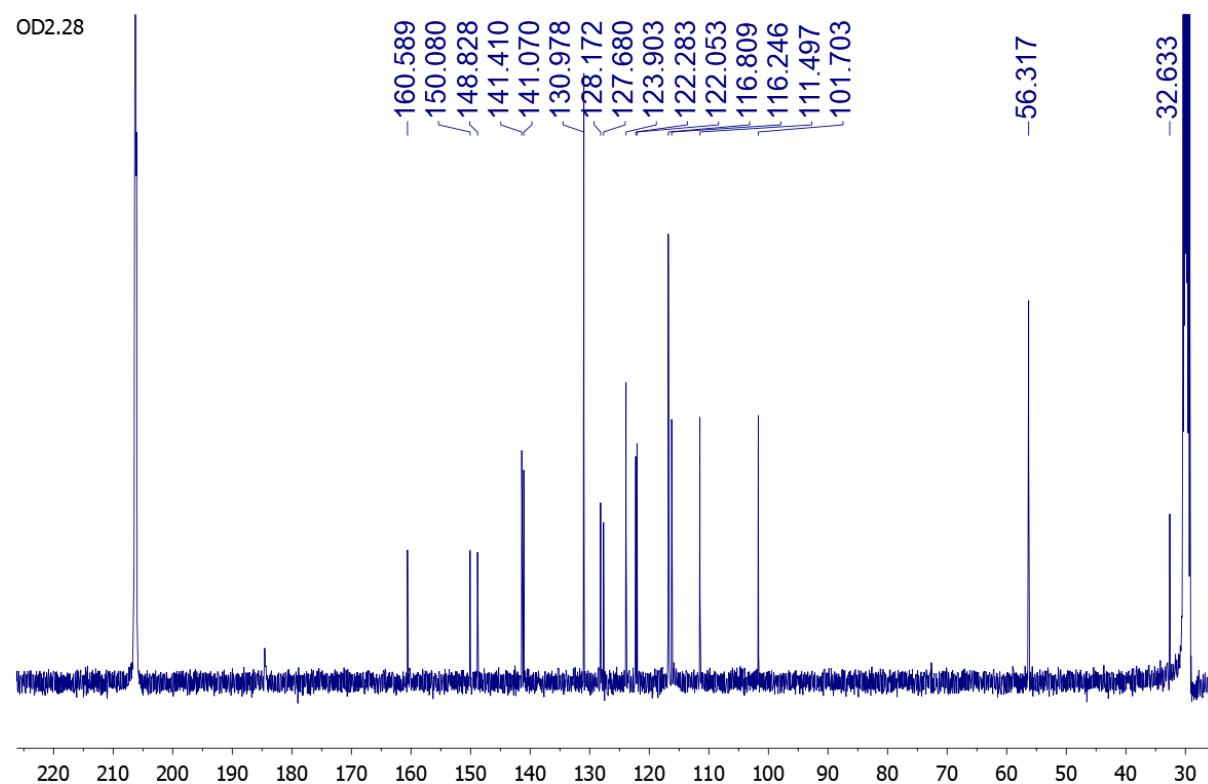
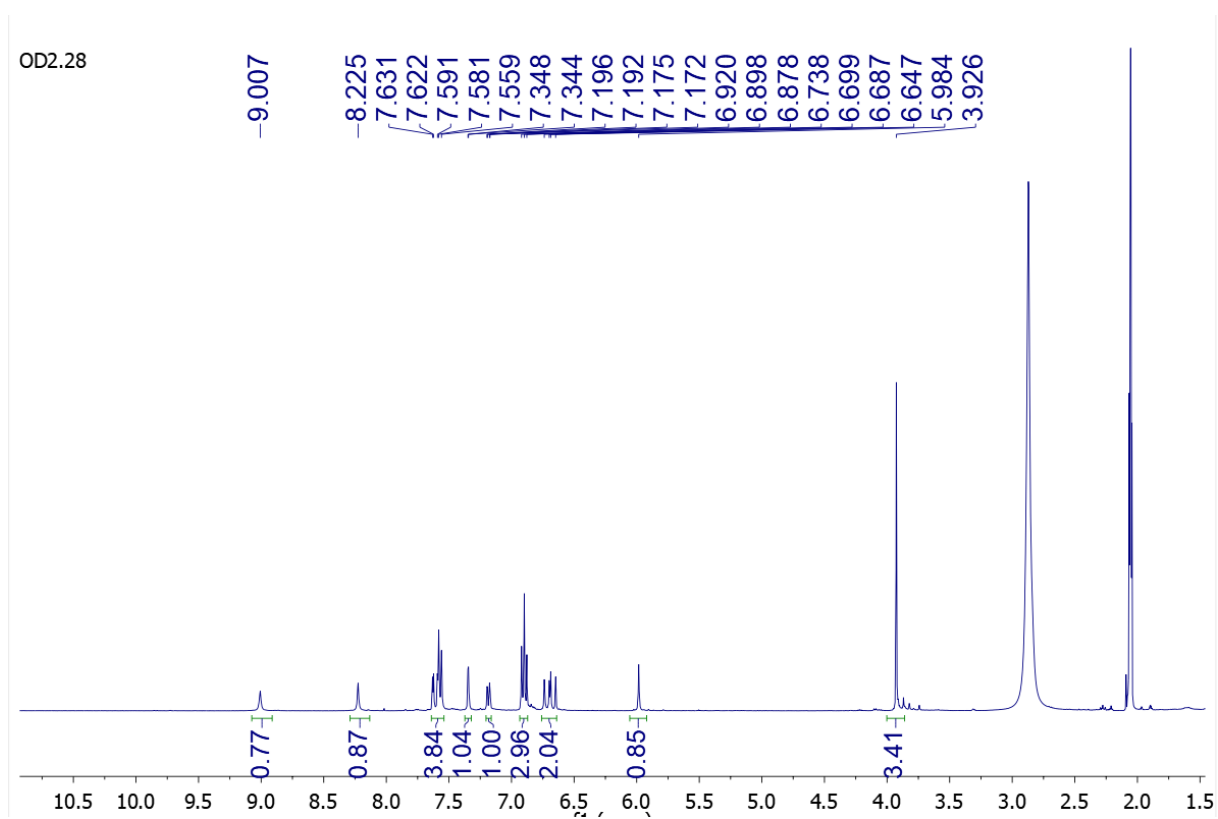


Figure S17: The ^1H and ^{13}C NMR spectrum of **5** in acetone- d_6

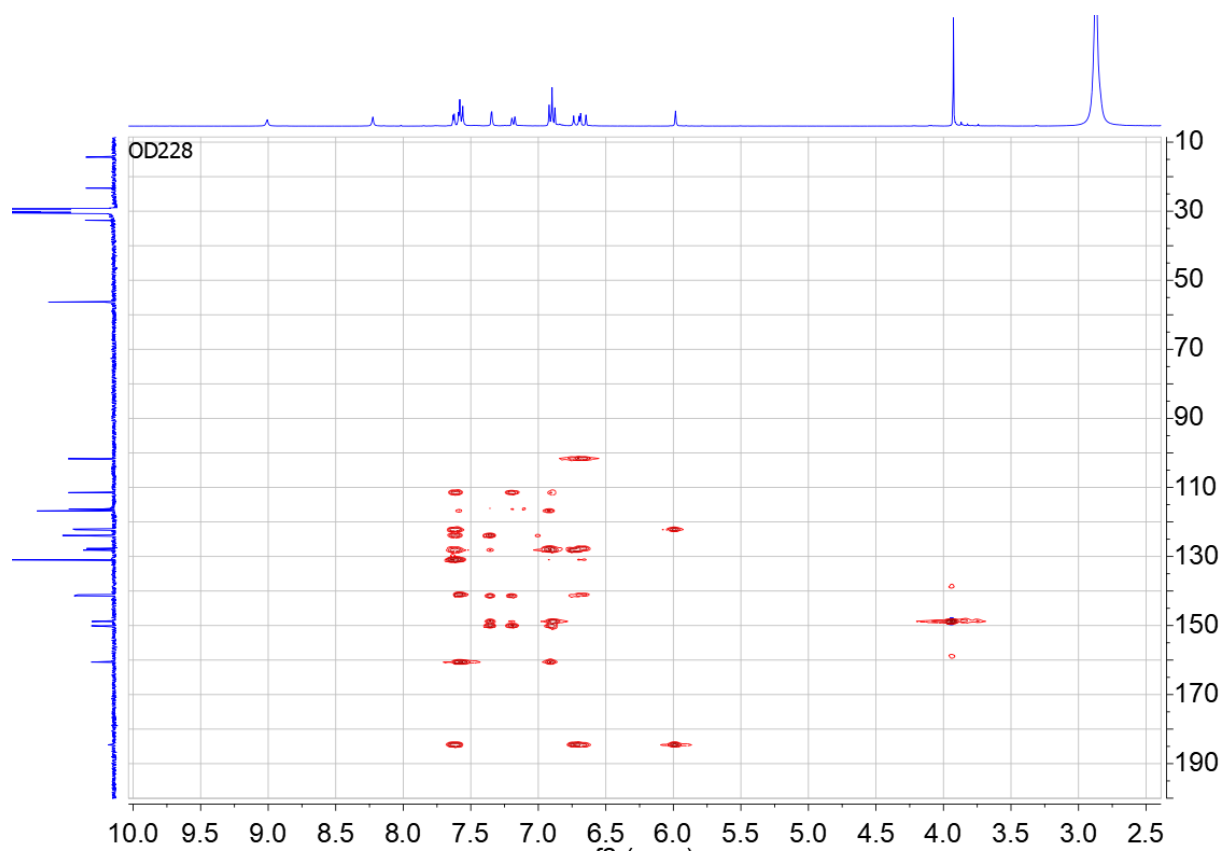


Figure S18: The HMBC spectrum of **5** in acetone- d_6

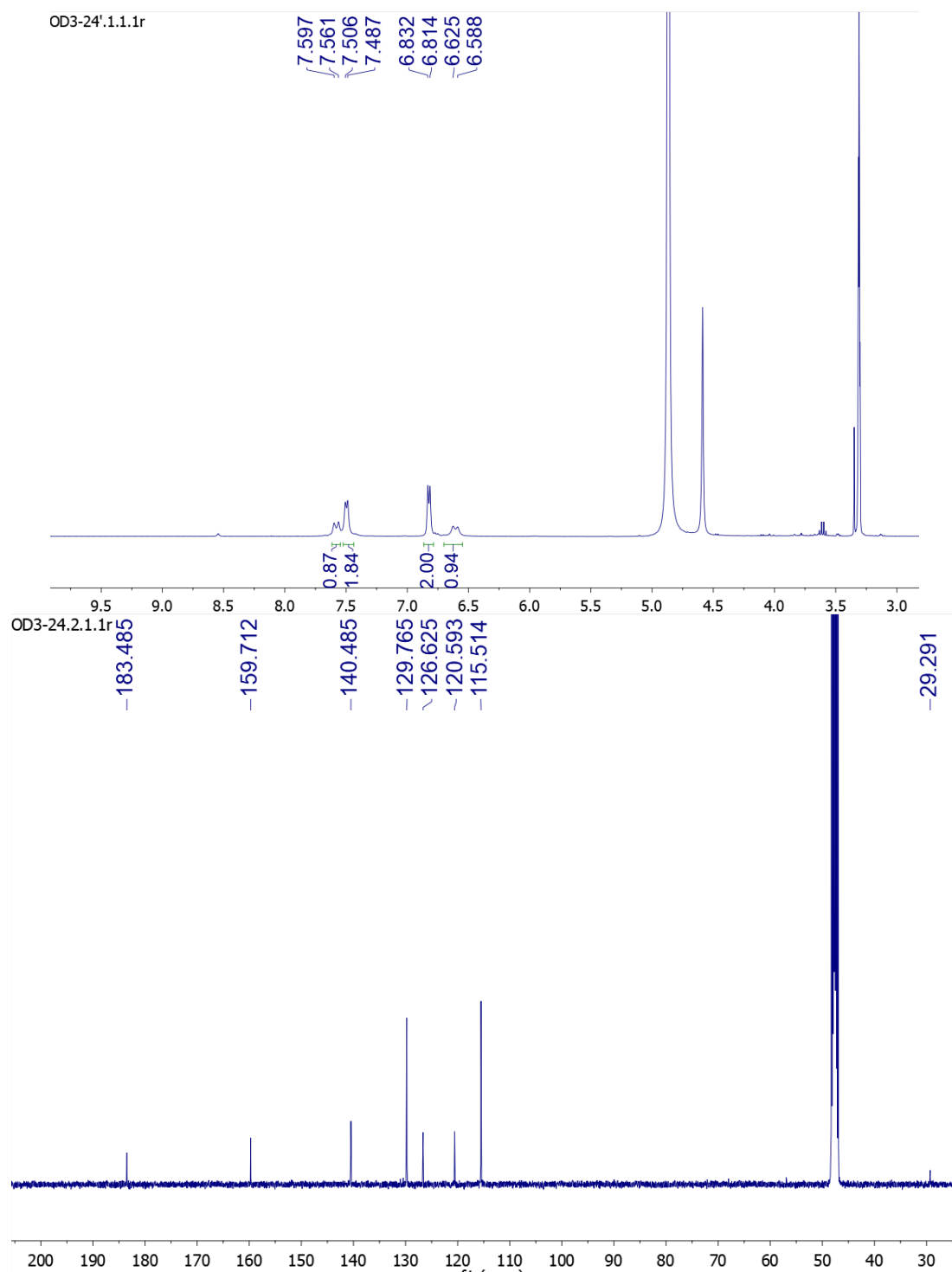


Figure S19: The ^1H and ^{13}C NMR spectrum of **6** in methanol- d_4

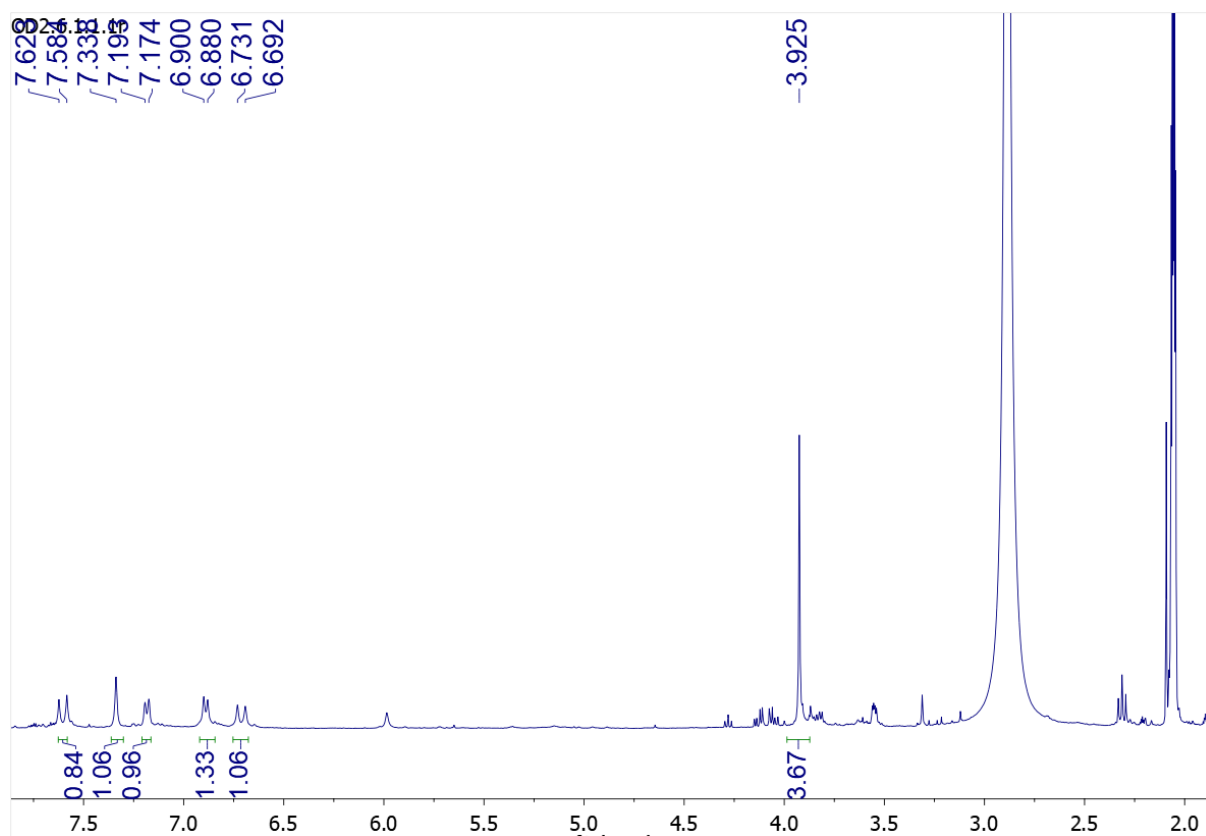


Figure S20: The ^1H NMR spectrum of **7** in acetone- d_6

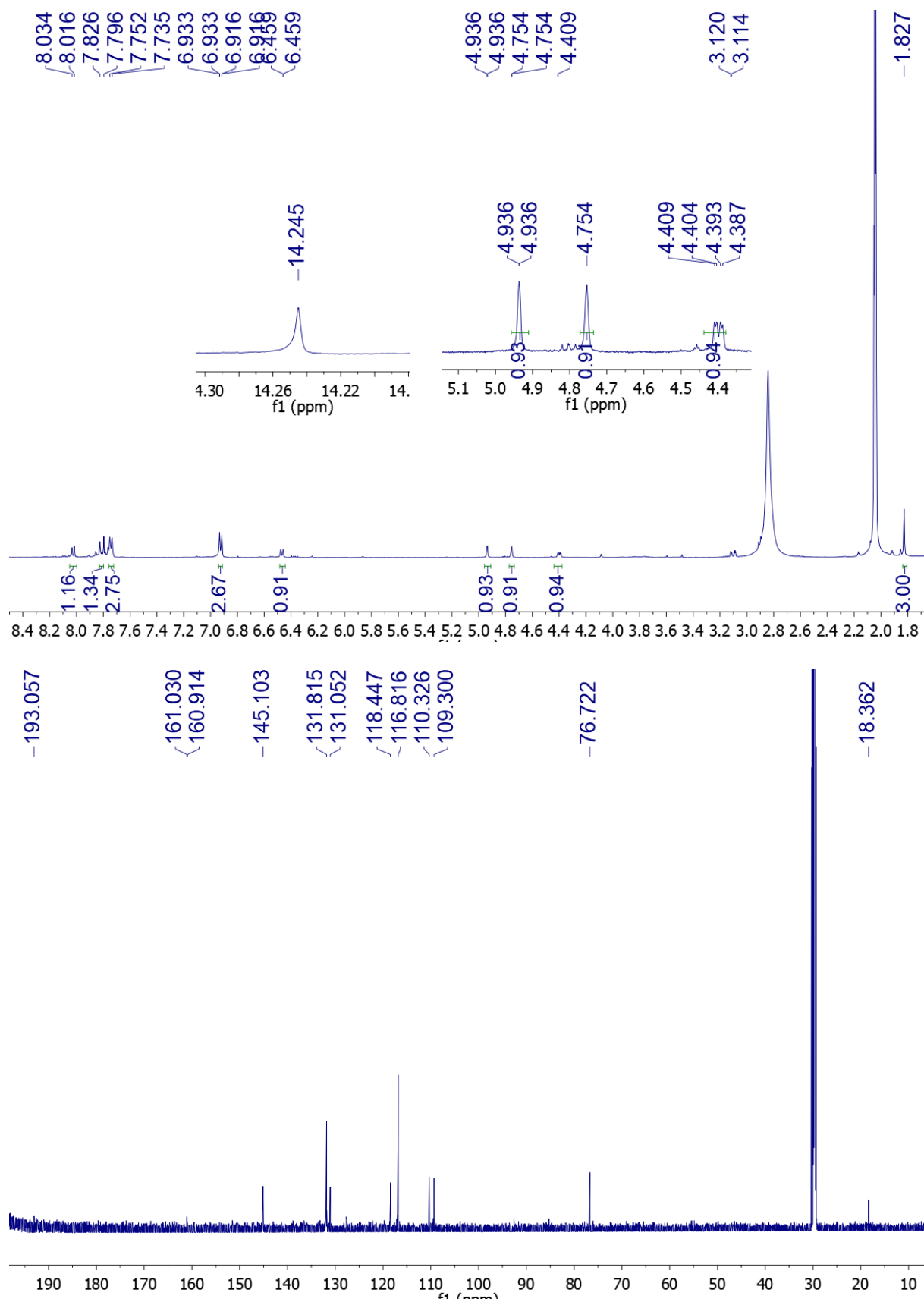


Figure S21: The ^1H and ^{13}C NMR spectrum of **8** in acetone- d_6

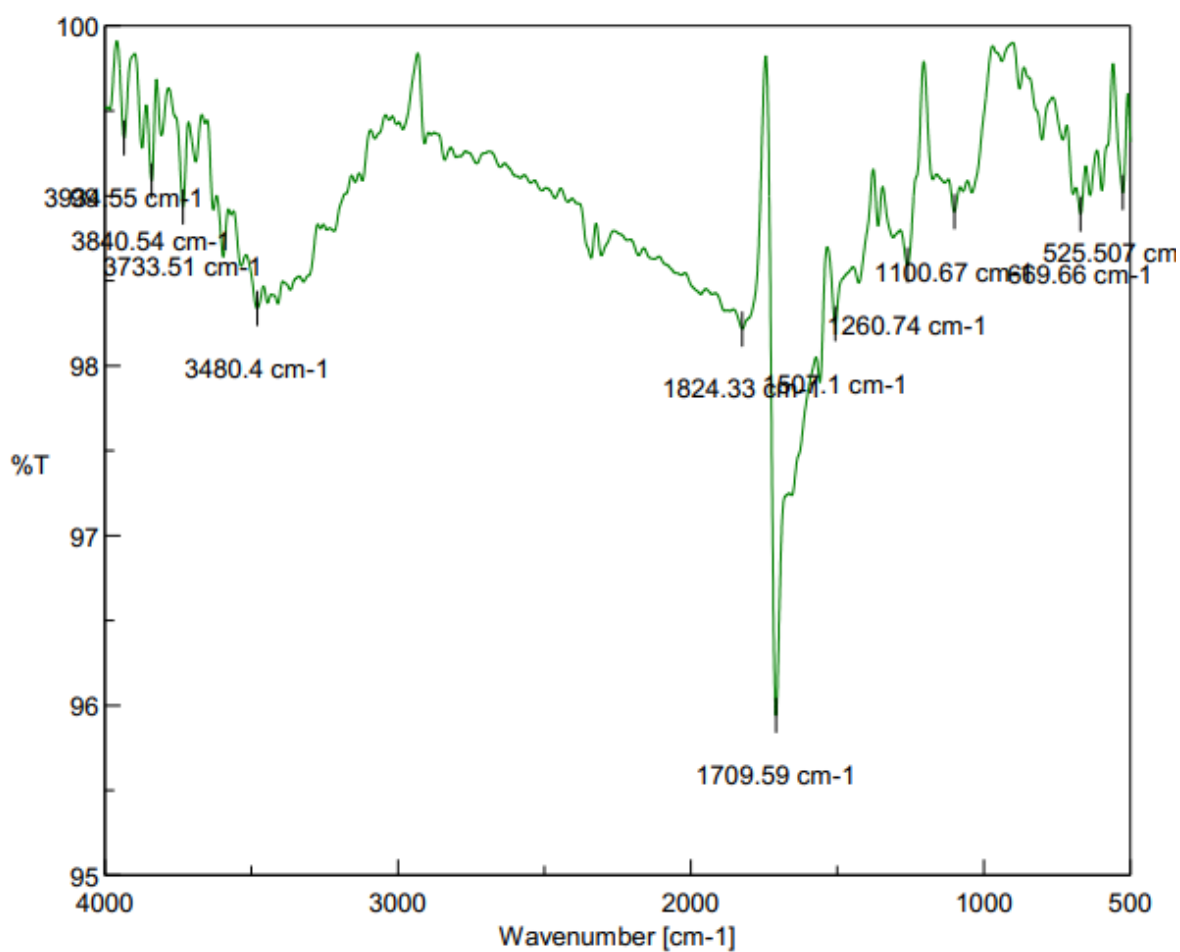
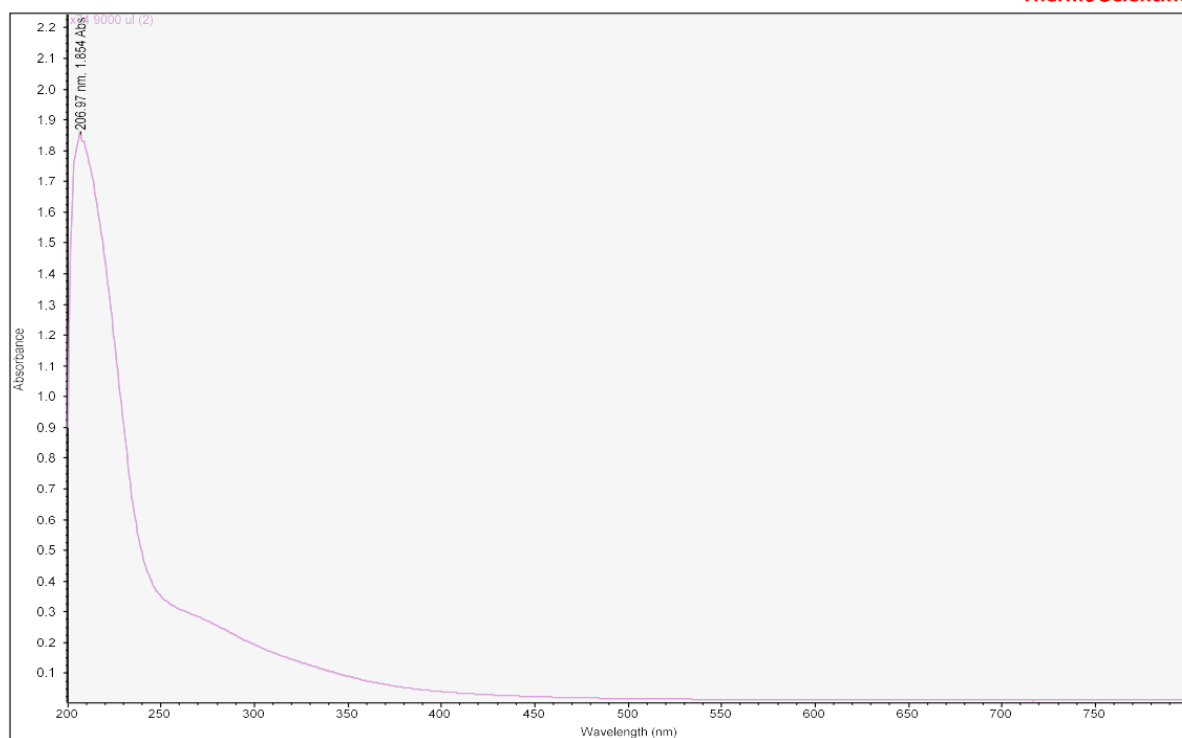


Figure S22: UV and IR spectra of **1**.

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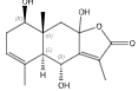
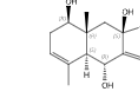
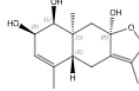
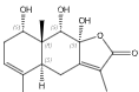
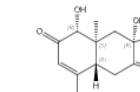
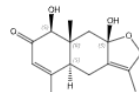
<p>1 96</p> <p>1174550-77-6</p>  <p>Available stereochemistry shown, Rotation (+)</p> <p>C₁₅H₂₀O₅ Linderagalactone E</p> <p>4 References 0 Reactions 1 Supplier</p>	<p>2 96</p> <p>2764604-96-6</p>  <p>Absolute stereochemistry shown, Rotation (+)</p> <p>C₁₅H₂₀O₅ Naphtho[2,3-b]furan-2(4H)-one, 4a,7,8,8a,9,9a-hexahydro-4,8,9a-trihydroxy-3,5,8a...</p> <p>1 Reference 0 Reactions 0 Suppliers</p>	<p>3 96</p> <p>2146087-86-5</p>  <p>Absolute stereochemistry shown, Rotation (-)</p> <p>C₁₅H₂₀O₅ Naphtho[2,3-b]furan-2(4H)-one, 4a,7,8,8a,9,9a-hexahydro-7,8,9a-trihydroxy-3,5,8a...</p> <p>1 Reference 0 Reactions 1 Supplier</p>
<p>4 95</p> <p>1384853-12-6</p>  <p>Absolute stereochemistry shown, Rotation (+)</p> <p>C₁₅H₂₀O₅ (4aS,8S,8aR,9S,9aS)-4a,7,8,8a,9,9a-hexahydro-8,9,9a-trihydroxy-3,5,8a-trimethyl...</p> <p>1 Reference 0 Reactions 0 Suppliers</p>	<p>5 95</p> <p>131740-68-6</p>  <p>Absolute stereochemistry shown</p> <p>C₁₅H₁₈O₅ Naphtho[2,3-b]furan-2,7-dione, 4,4a,8,8a,9,9a-hexahydro-8,9a-dihydroxy-3,5,8a-tr...</p> <p>3 References 2 Reactions 0 Suppliers</p>	<p>6 95</p> <p>1664367-04-7</p>  <p>Absolute stereochemistry shown, Rotation (-)</p> <p>C₁₅H₁₈O₅ (4aS,8S,8aR,9aS)-4,4a,8,8a,9,9a-hexahydro-8,9a-dihydroxy-3,5,8a-trimethylnaphtho...</p> <p>1 Reference 0 Reactions 1 Supplier</p>

Figure S23: Scifinder searching for 1.

Table S1. ^1H NMR (500 MHz, δ_{H} , multi, (J in Hz) and ^{13}C NMR (125 MHz) spectral data of comparison of compound **1** and eudebeiolide J

No.	1 (DMSO- d_6)		Eudebeiolide J [1] (methanol- d_4)	
	δ_{H} (J in Hz)	δ_{C}	δ_{H} (J in Hz)	δ_{C}
1	3.21, d, $J = 4.0$	76.6	3.37, d, $J = 4.2$	75.8
2	3.84, m	70.4	4.24, brs	68.0
3	5.57, brs	122.9	5.31, s	124.9
4		137.7		137.1
5	2.21, d, $J = 13.5$	39.9	2.36, brd, $J = 13.5$	43.8
6	2.96, d, $J = 10.0$ 2.18, dd, $J = 13.5, 10.0$	22.3	2.95, dd, $J = 13.2, 3.6$ 2.26, td, $J = 13.2, 1.2$	24.3
7		161.6		163.0
8		104.6		106.1
9	1.88, d, $J = 13.5$ 1.86, d, $J = 13.5$	45.4	2.07, d, $J = 13.2$ 1.97, d, $J = 13.2$	46.1
10		36.8		42.0
11		120.4		122.9
12		171.8		174.6
13	1.73, s	7.9	1.83, s	8.2
14	1.03, s	16.1	1.13, s	16.4
15	4.01, dd, $J = 13.5, 5.5$ 3.90, dd, $J = 13.5, 5.5$	62.8	1.78, d, $J = 0.6$	21.4
1-OH	4.85, d, $J = 4.0$			
2-OH	4.80, d, $J = 4.5$			
8-OH	7.12, s			
15-OH	4.75, t, $J = 5.5$			

[1] H.-J. Jang, S. Lee, S.-J. Lee, H.-J. Lim, K. Jung, Y.H. Kim, S.W. Lee and M.-C. Rho (2017). Anti-inflammatory activity of eudesmane-type sesquiterpenoids from *Salvia plebeia*, *J. Nat. Prod.* 80, 2666-2676.