

## Supporting Information

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### A New Dimeric Sesquiterpenoid from *Chloranthus japonicus* Sieb.

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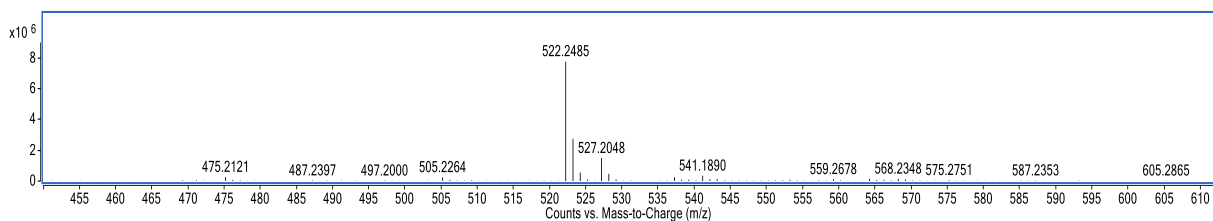
\* Corresponding author: E-Mail: hope1126@hotmail.com; Phone: 086-21-57423899

Qi-Yan Li and Yan Wang contributed equally to this work.

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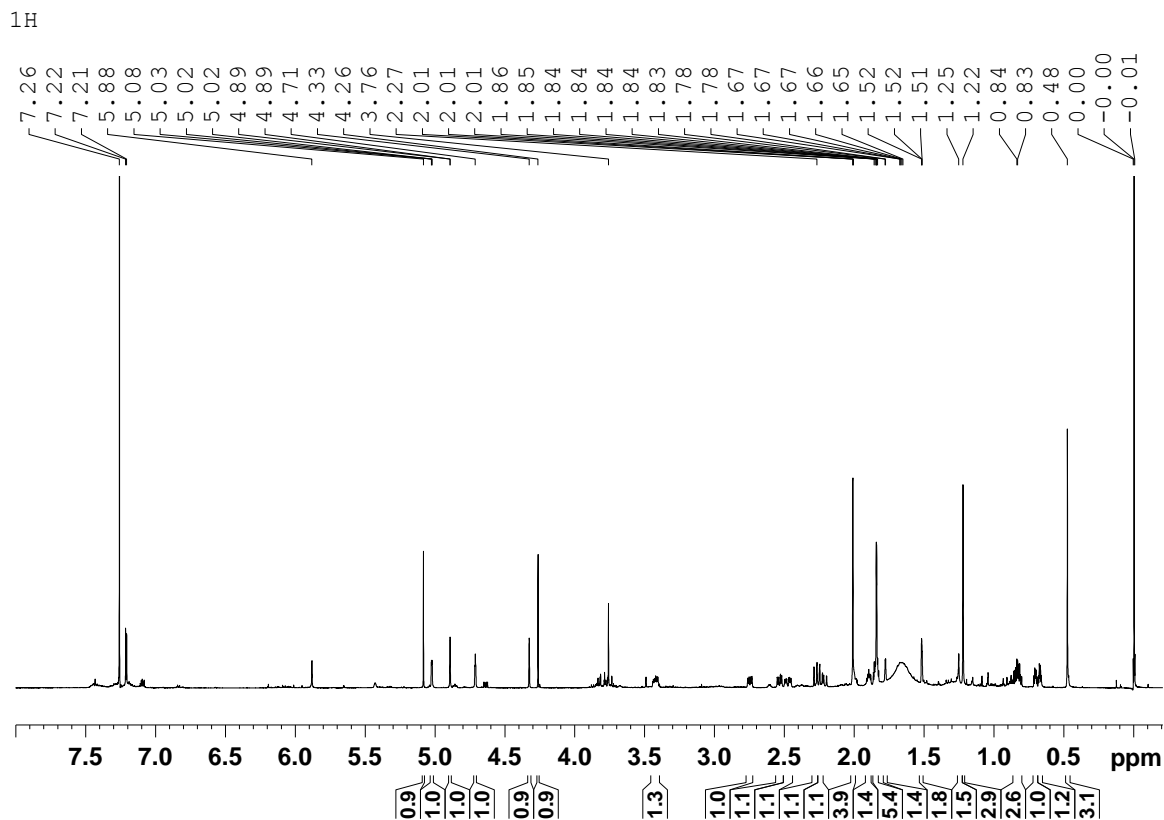
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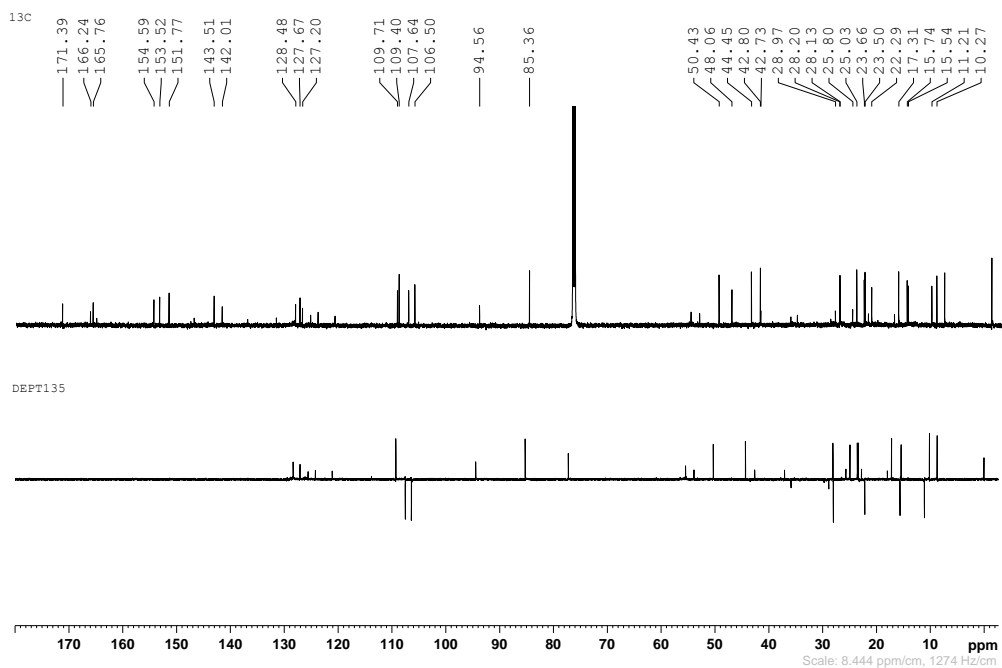


Formula	Score (MFG)	Mass	Mass (MFG)	$m/z$ (calcd.)	Diff (ppm)	$m/z$
$C_{30}H_{32}O_7$	100	504.2147	504.2148	522.2486	0.26	522.2485
$C_{30}H_{32}O_7$	99.96	504.2156	504.2148	527.2040	-1.54	527.2048

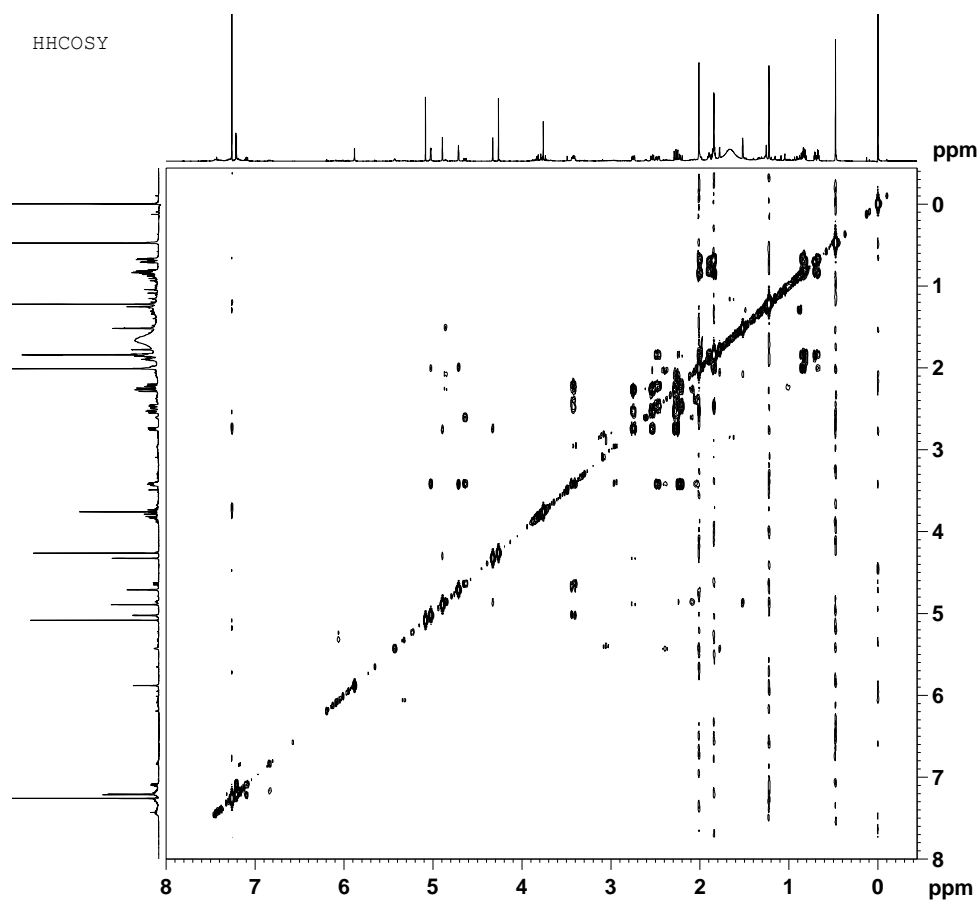
**Figure S1:** HRESIMS spectrum of **1**



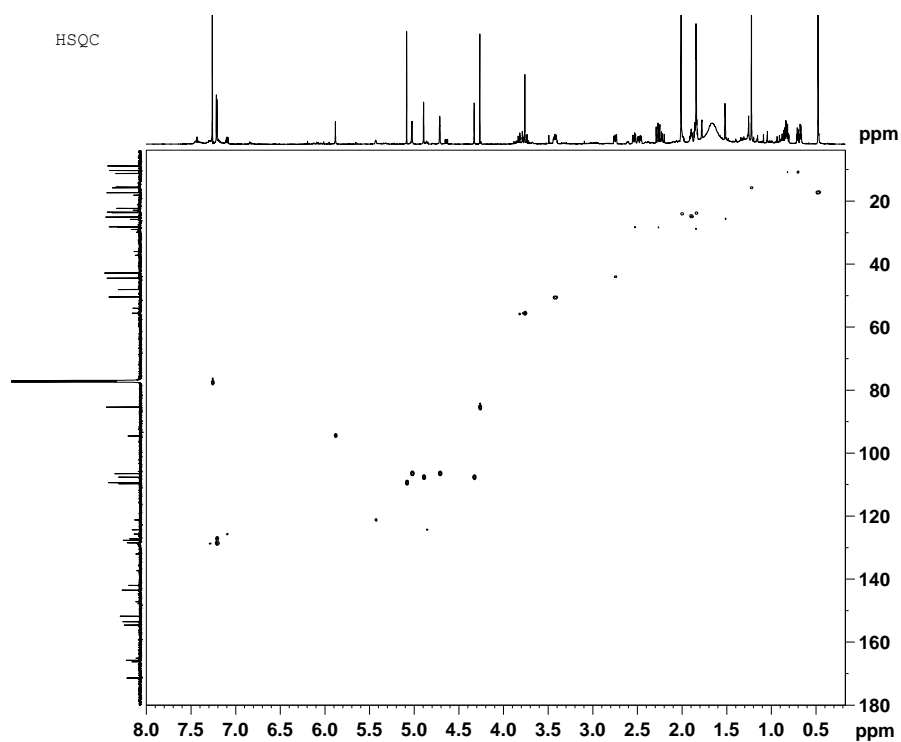
**Figure S2:**  $^1H$  NMR spectrum of **1**



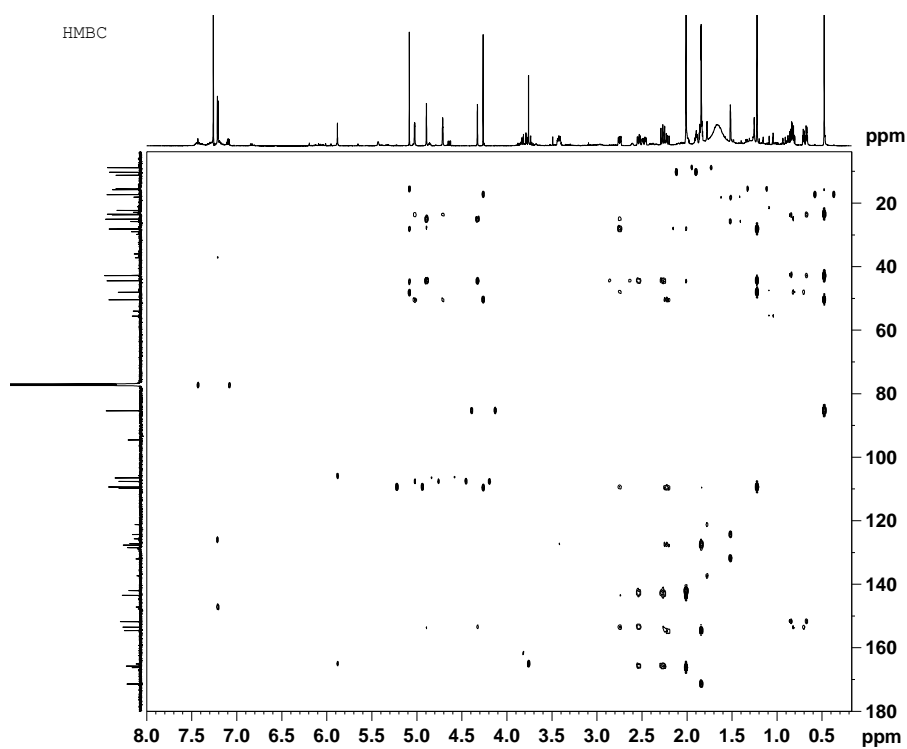
**Figure S3:** <sup>13</sup>C NMR and DEPT135 spectra of **1**



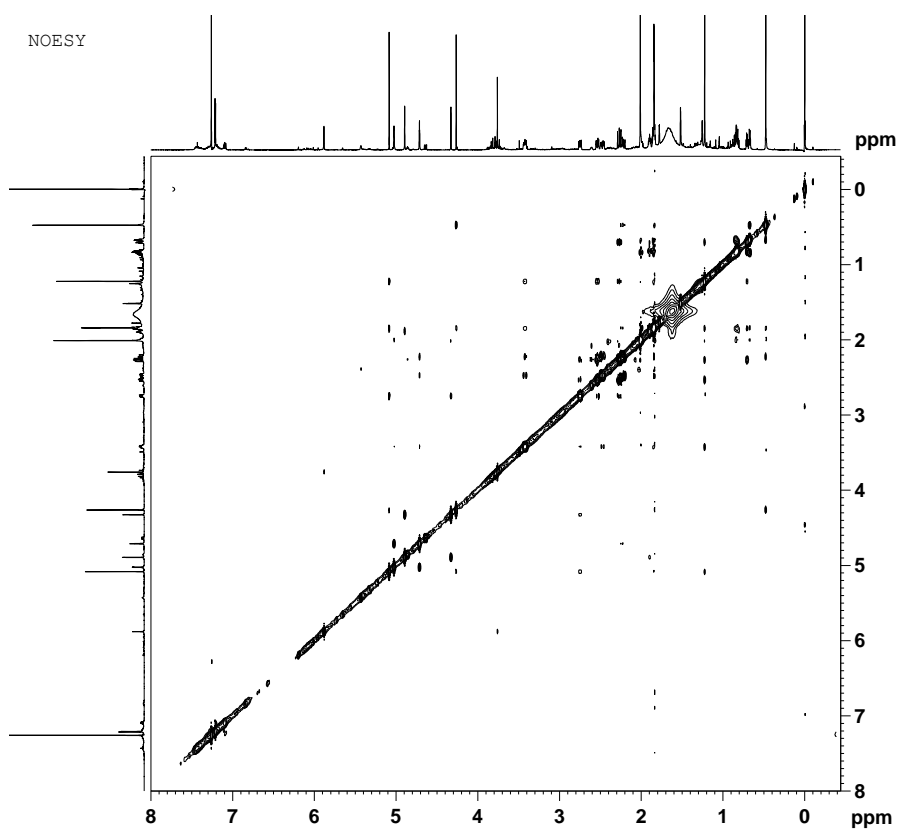
**Figure S4:** <sup>1</sup>H-<sup>1</sup>H COSY spectrum of **1**



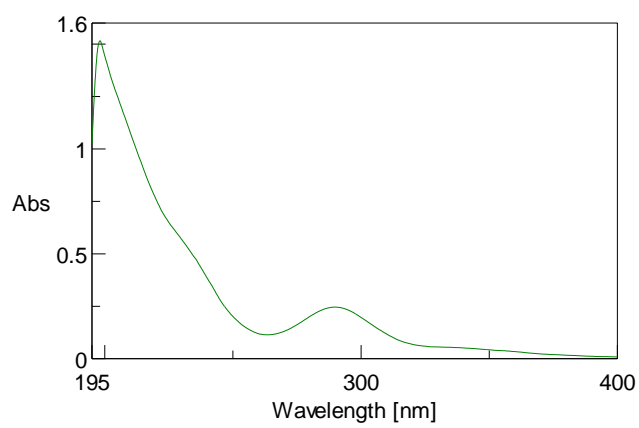
**Figure S5: HSQC spectrum of 1**



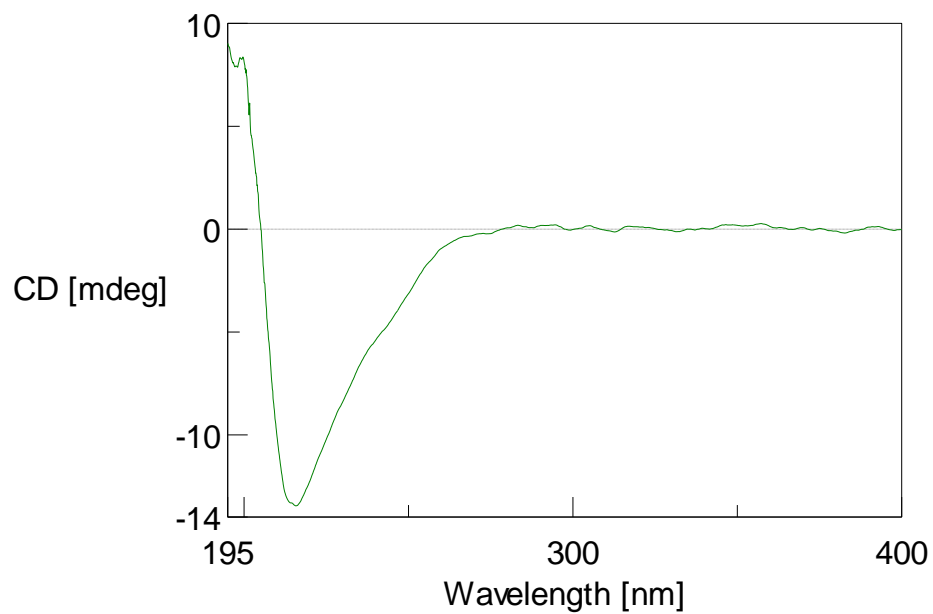
**Figure S6: HMBC spectrum of 1**



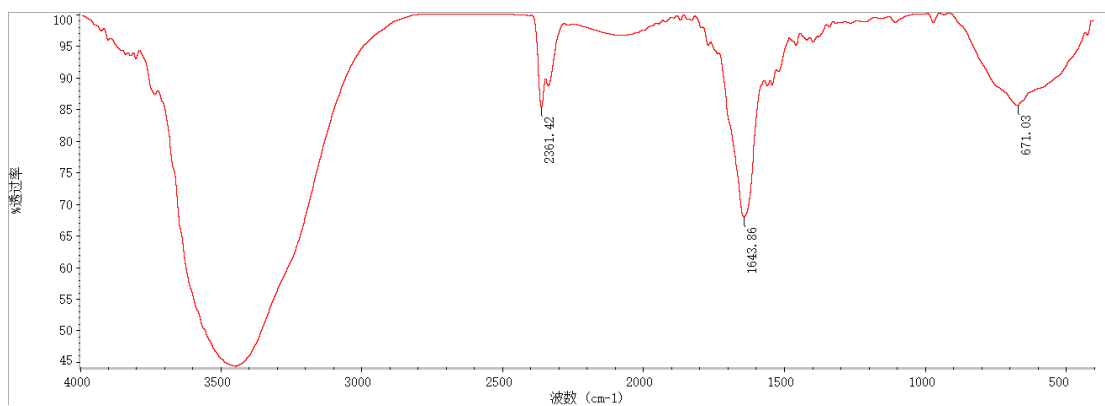
**Figure S7:** NOESY spectrum of **1**



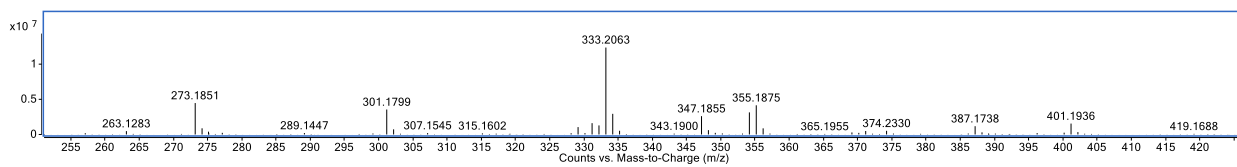
**Figure S8:** UV spectrum of **1**



**Figure S9: ECD spectrum of 1**



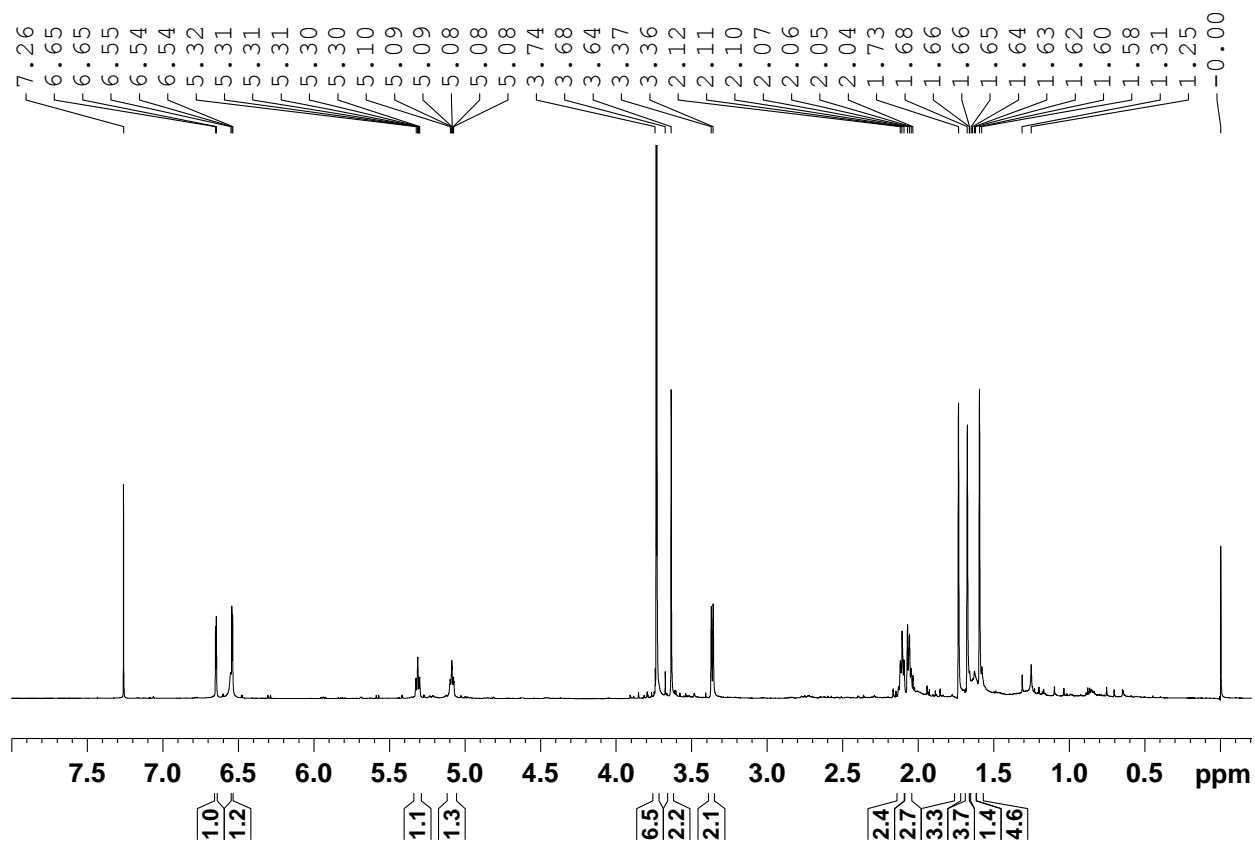
**Figure S10: IR spectrum of 1**



Formula	Score (MFG)	Mass	Mass (MFG)	$m/z$ (calcd.)	Diff (ppm)	$m/z$
C <sub>20</sub> H <sub>28</sub> O <sub>4</sub>	99.99	332.1990	332.1988	333.2060	-0.80	333.2063
C <sub>20</sub> H <sub>28</sub> O <sub>4</sub>	99.97	332.1983	332.1988	355.1880	1.45	355.1875

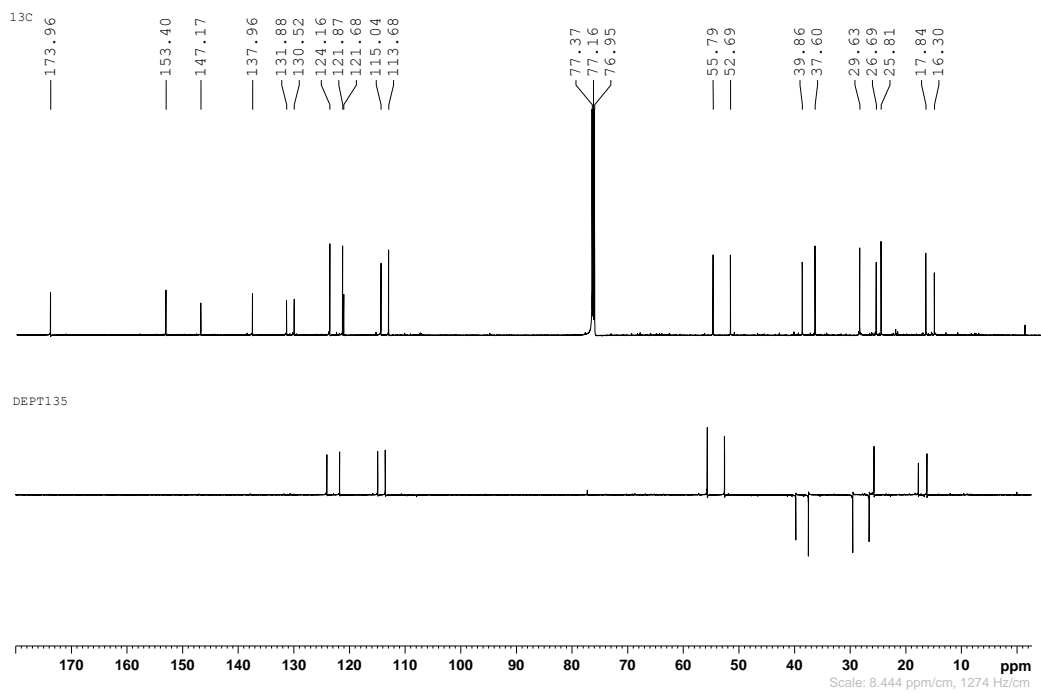
**Figure S11: HRESIMS spectrum of 2**

<sup>1</sup>H

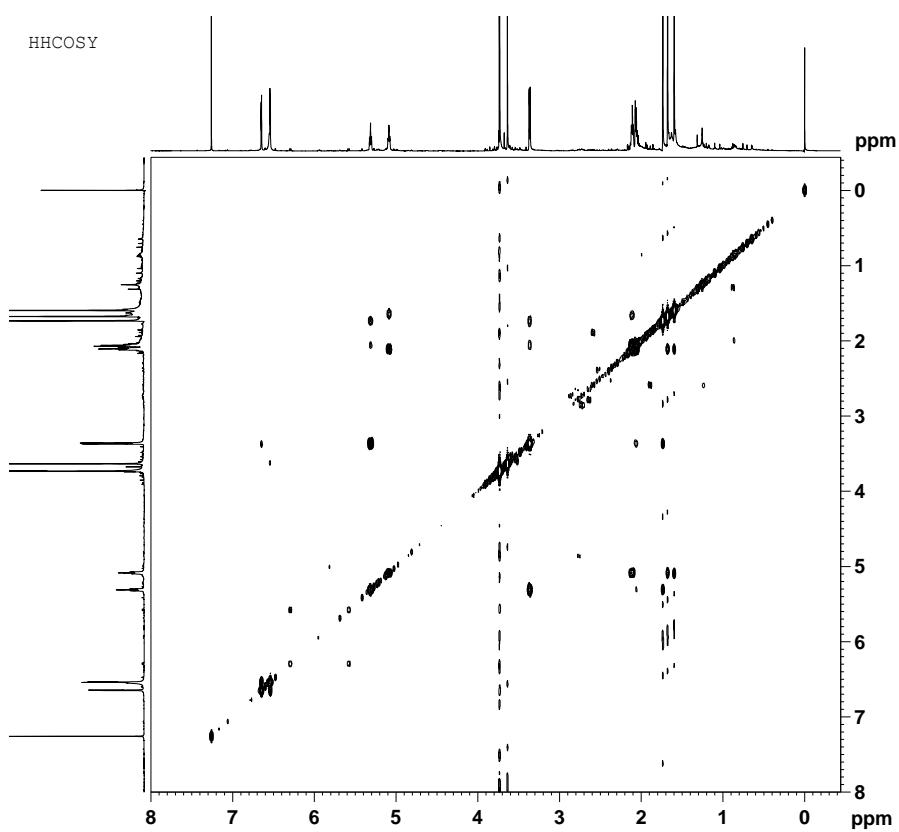


**Figure S12: <sup>1</sup>H NMR spectrum of 2**

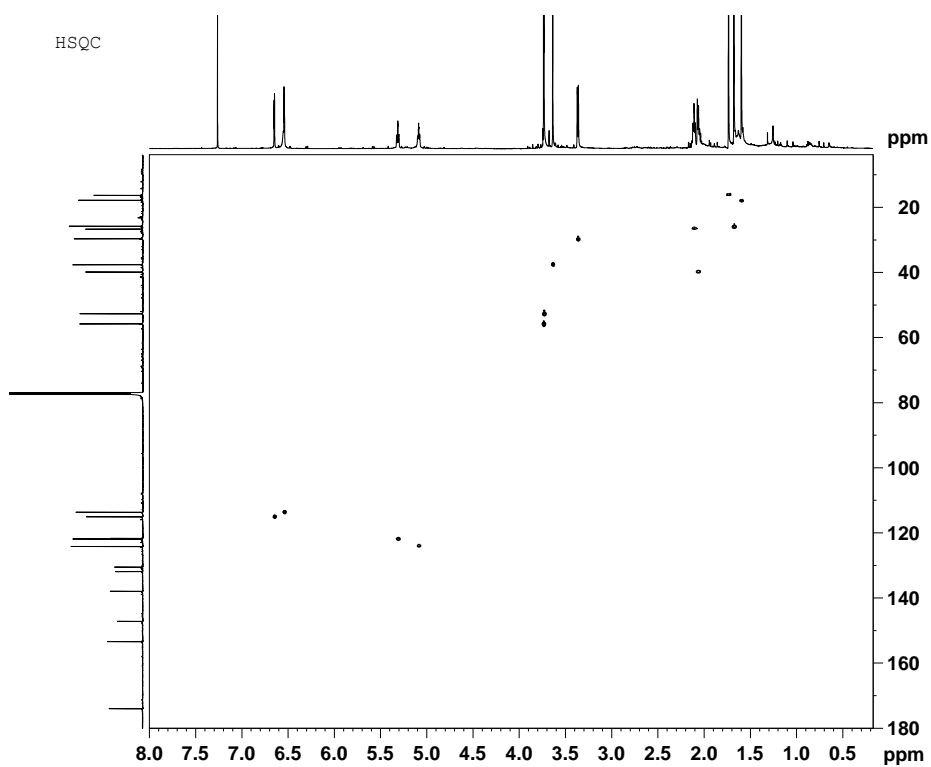




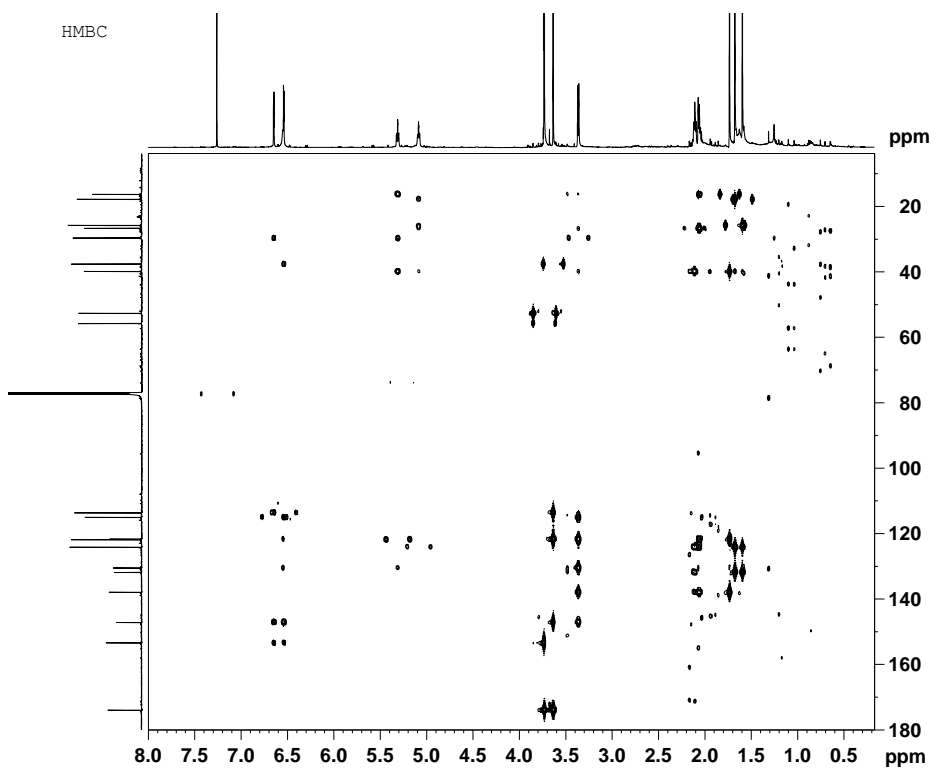
**Figure S13:** <sup>13</sup>C NMR and DEPT135 spectra of **2**



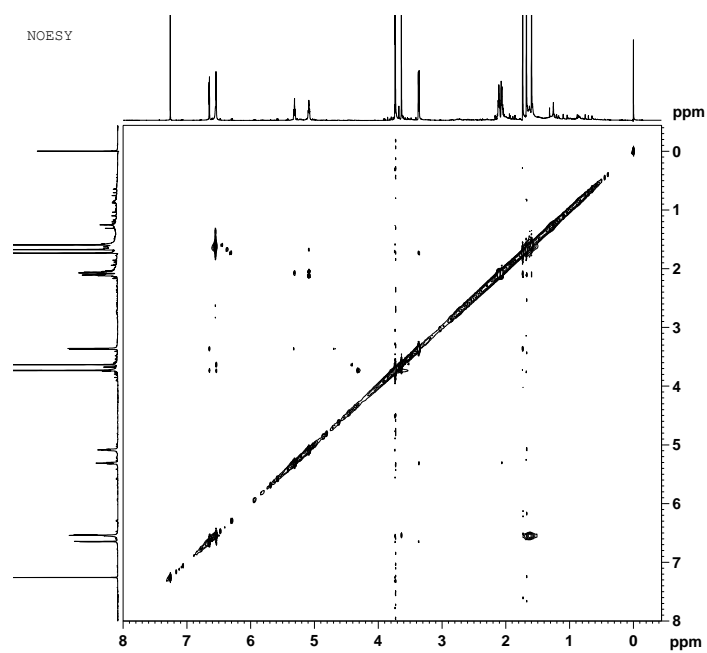
**Figure S14:** <sup>1</sup>H-<sup>1</sup>H COSY spectrum of **2**



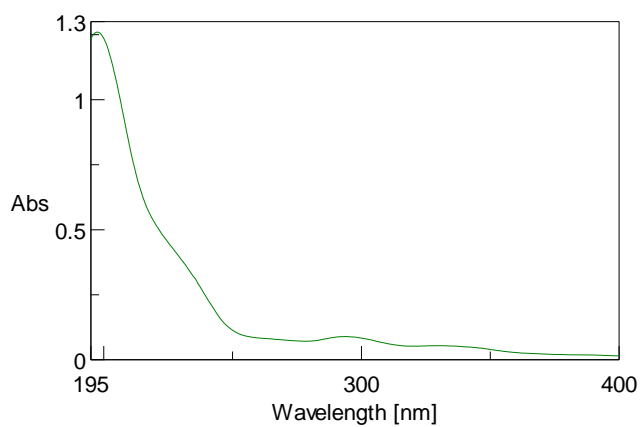
**Figure S15:** HSQC spectrum of **2**



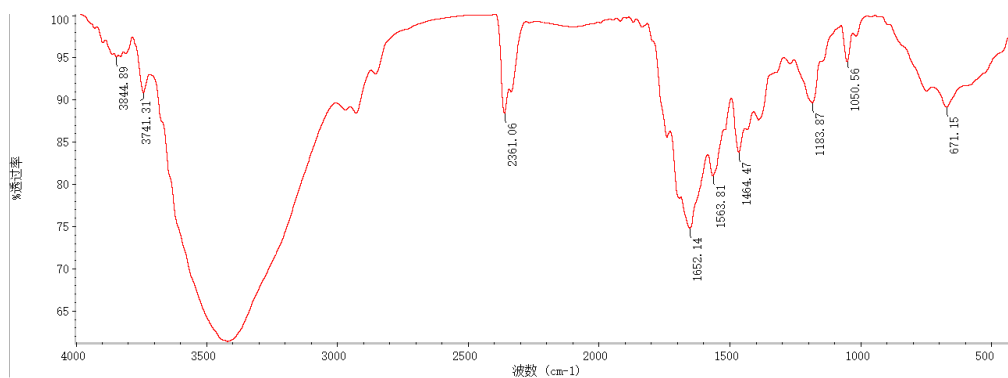
**Figure S16:** HMBC spectrum of **2**



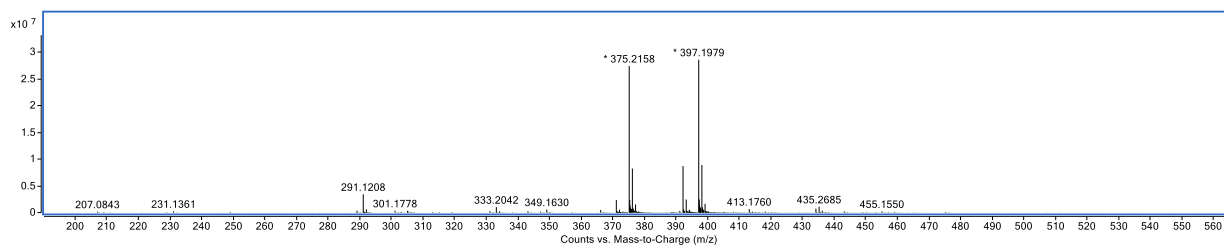
**Figure S17:** NOESY spectrum of **2**



**Figure S18:** UV spectrum of **2**



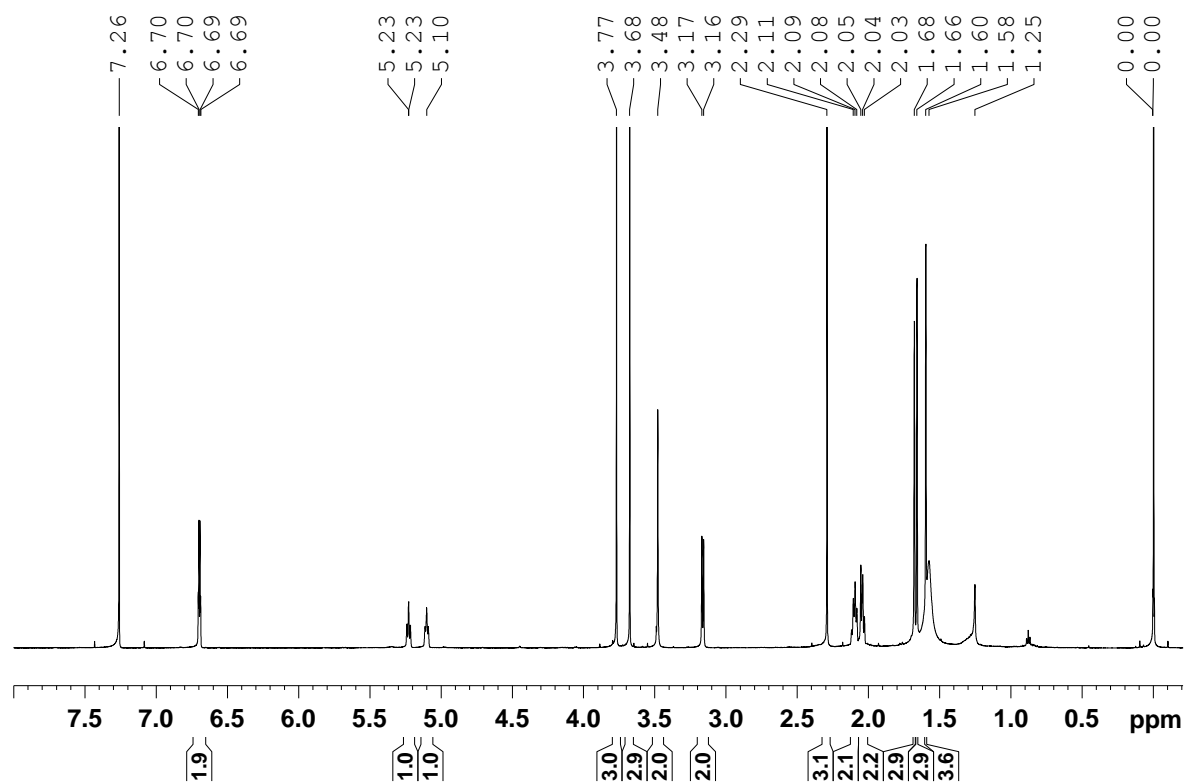
**Figure S19:** IR spectrum of **2**



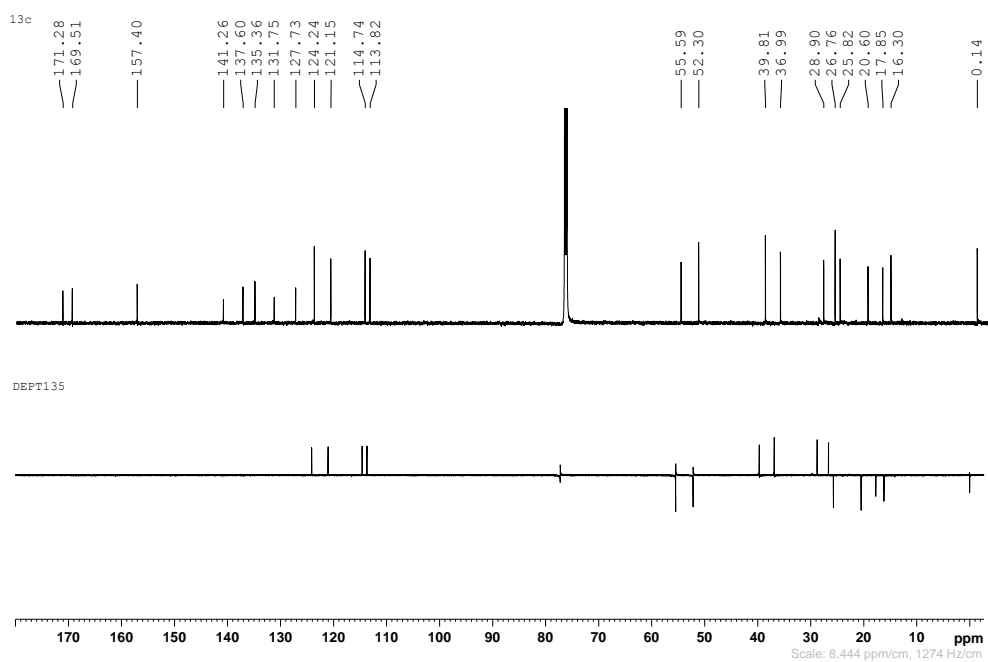
Formula	Score (MFG)	Mass	Mass (MFG)	$m/z$ (calcd.)	Diff (ppm)	$m/z$
C <sub>22</sub> H <sub>30</sub> O <sub>5</sub>	99.93	374.2085	374.2093	375.2166	2.14	375.2158
C <sub>22</sub> H <sub>30</sub> O <sub>5</sub>	99.95	374.2087	374.2093	397.1985	1.72	397.1979

**Figure S20:** HRESIMS spectrum of **3**

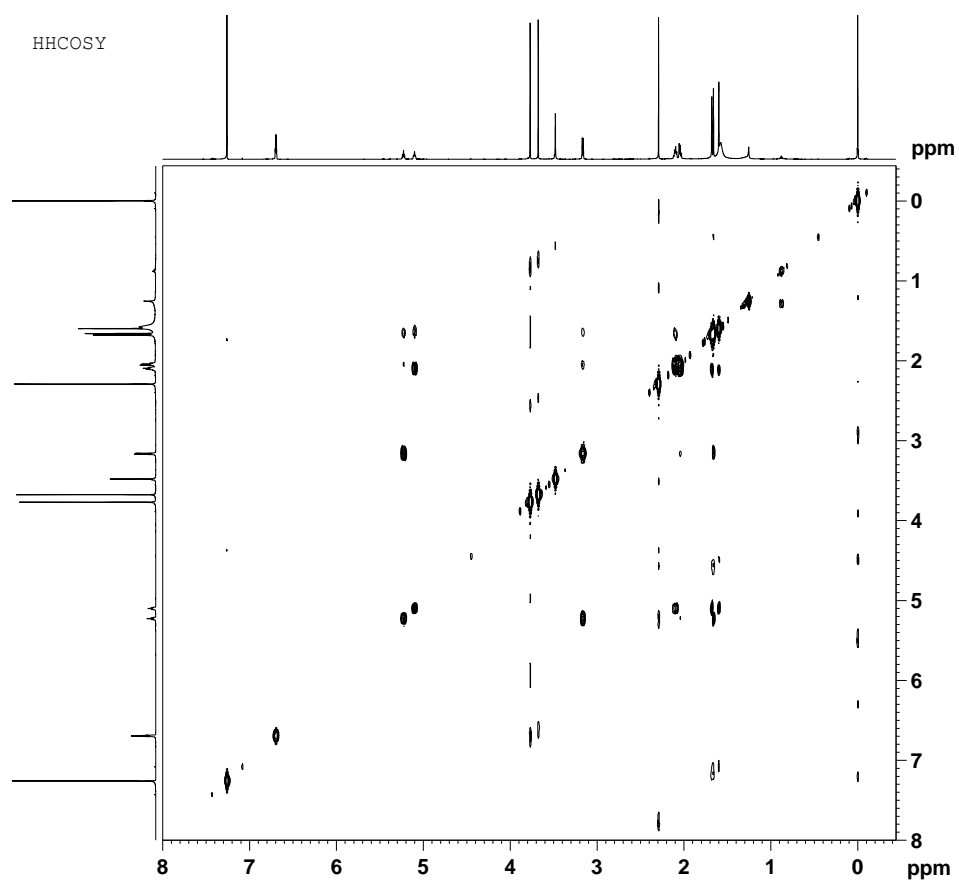
<sup>1</sup>H



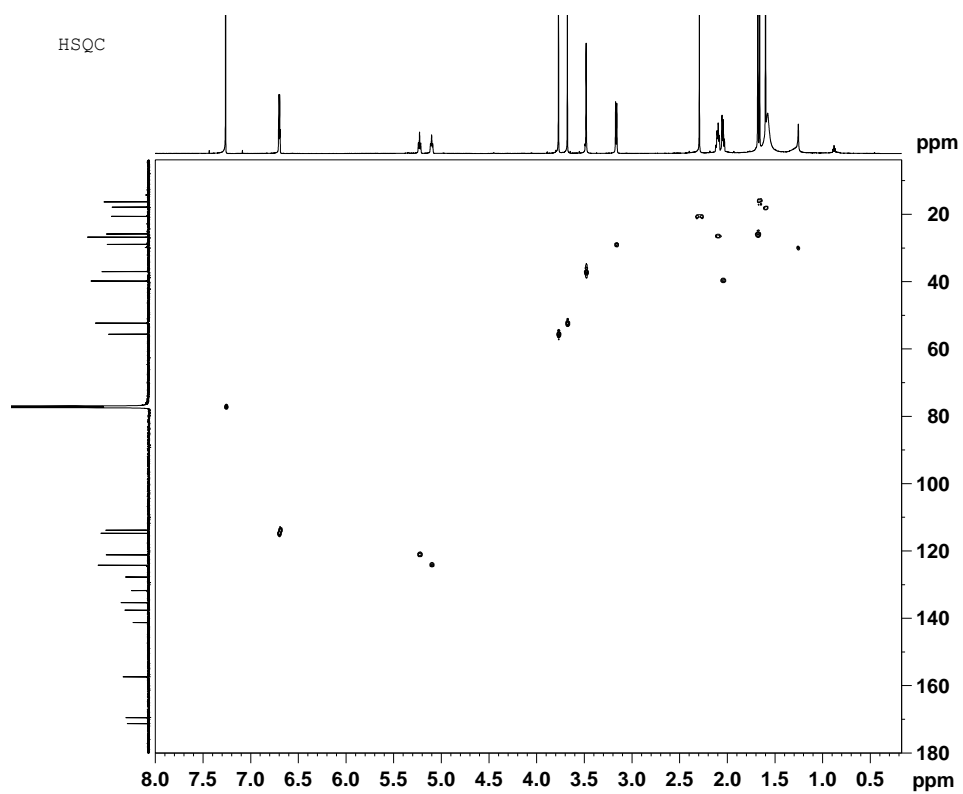
**Figure S21:** <sup>1</sup>H NMR spectrum of **3**



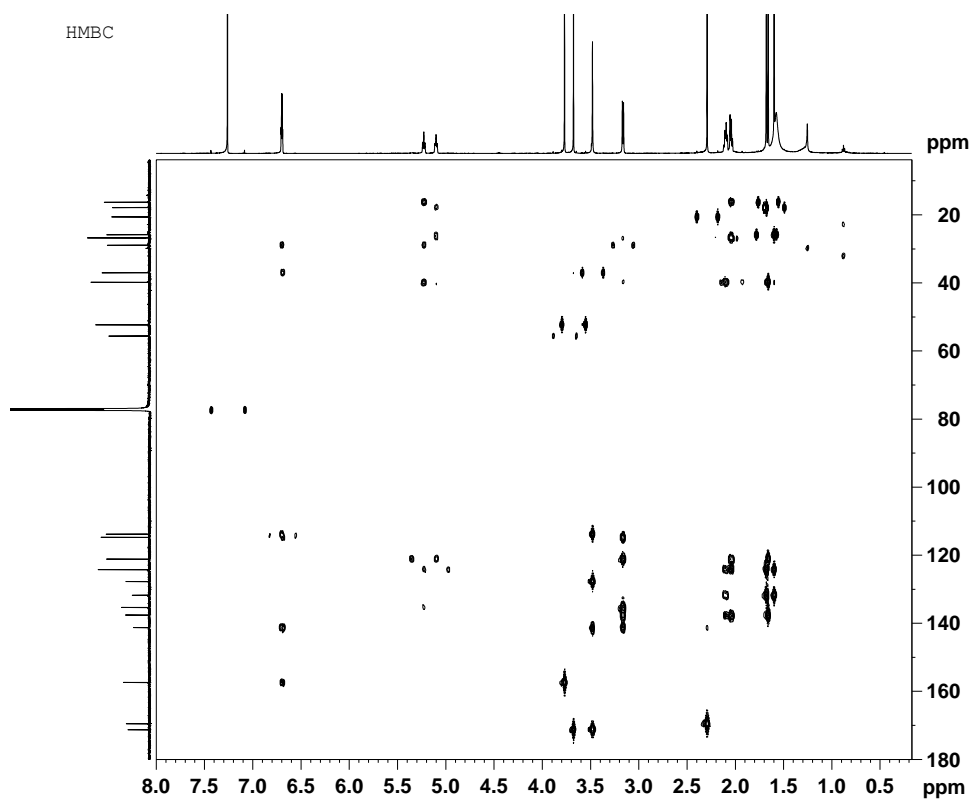
**Figure S22:** <sup>13</sup>C NMR and DEPT135 spectra of **3**



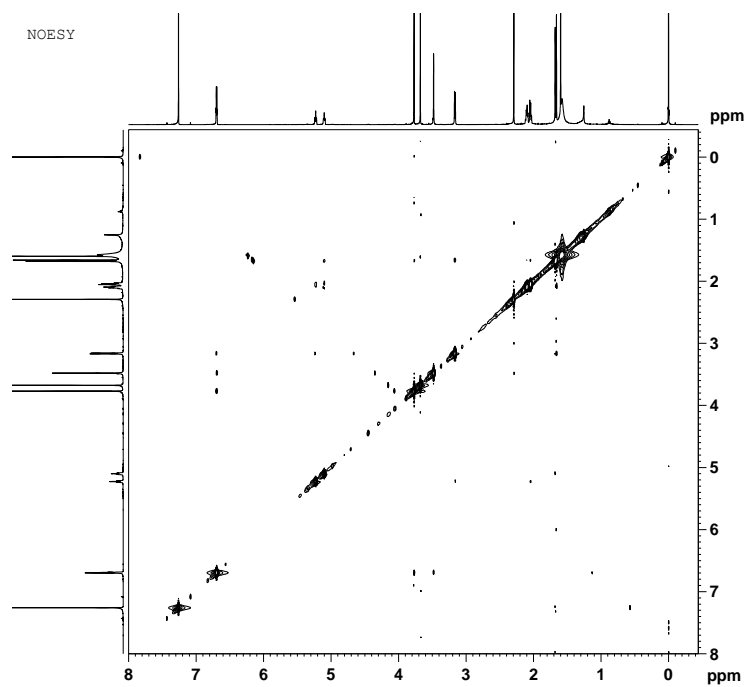
**Figure S23:** <sup>1</sup>H-<sup>1</sup>H COSY spectrum of **3**



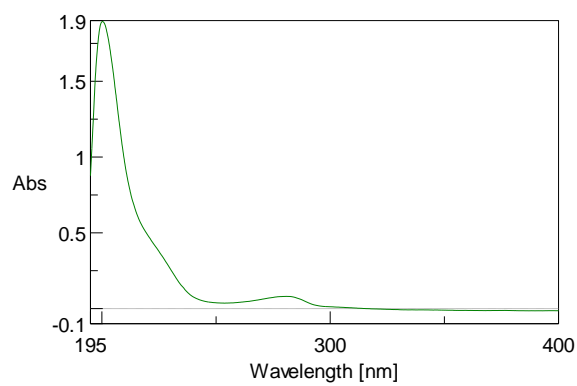
**Figure S24:** HSQC spectrum of **3**



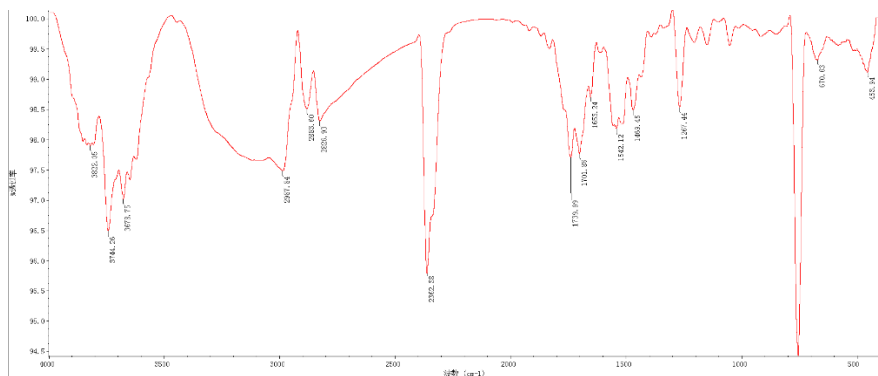
**Figure S25:** HMBC spectrum of **3**



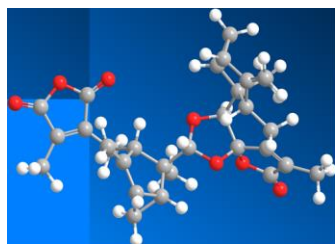
**Figure S26:** NOESY spectrum of **3**



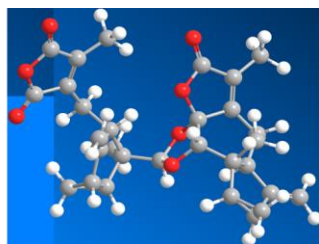
**Figure S27:** UV spectrum of **3**



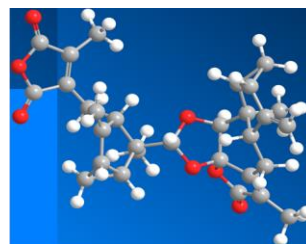
**Figure S28:** IR spectrum of **3**



**1-1 (50.15%)**



**1-2 (37.97%)**



**1-1 (11.88%)**

**Figure S29:** Optimized conformers (>1%) of **1**