

Supporting Information

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Two New Alkaloids from *Pleurotus ostreatus* (Jacq. :Pers.) Roll

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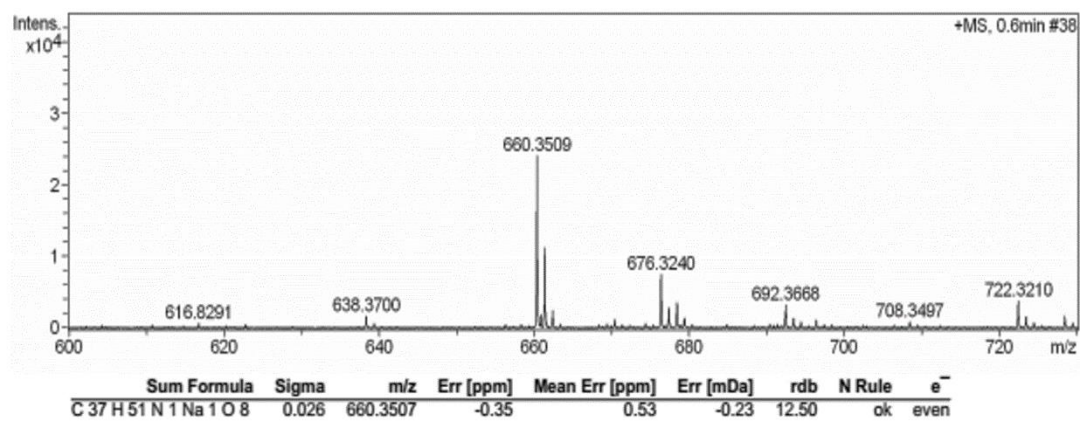


Figure S1: HR-ESI MS spectrum of **1** (Terpendole N)

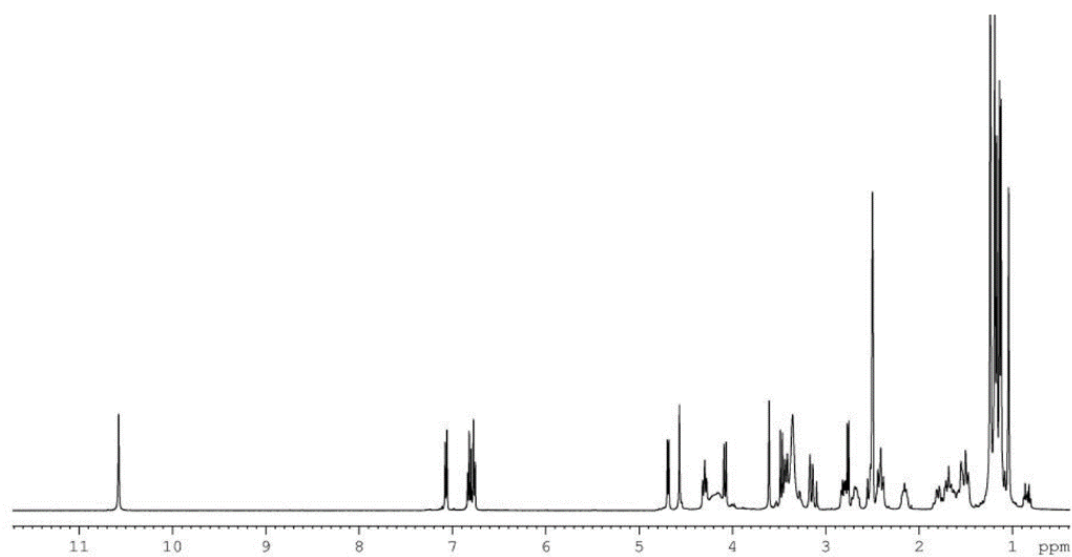


Figure S2: ^1H NMR (400 MHz, $\text{DMSO}-d_6$) spectrum of **1** (Terpendole N)

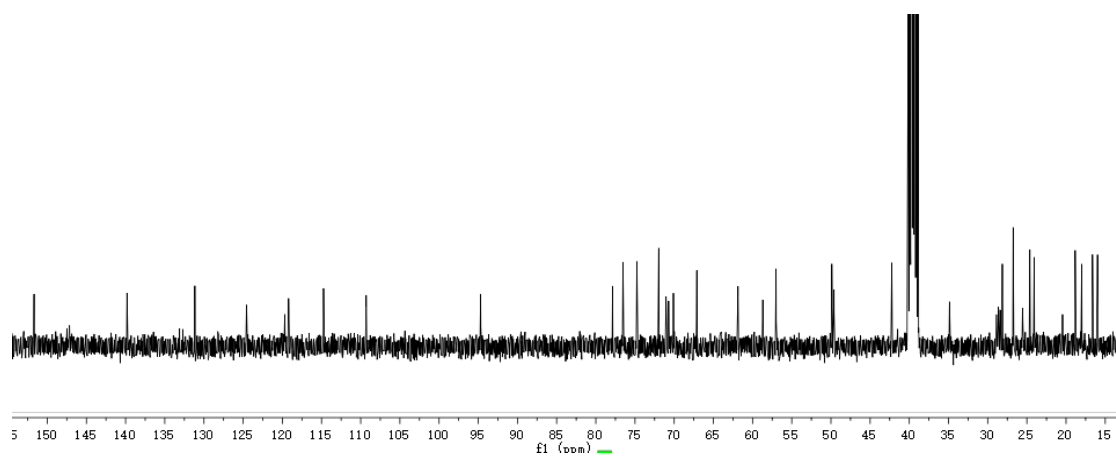


Figure S3: ^{13}C -NMR (100 MHz, $\text{DMSO-}d_6$) spectrum of **1** (Terpendole N)

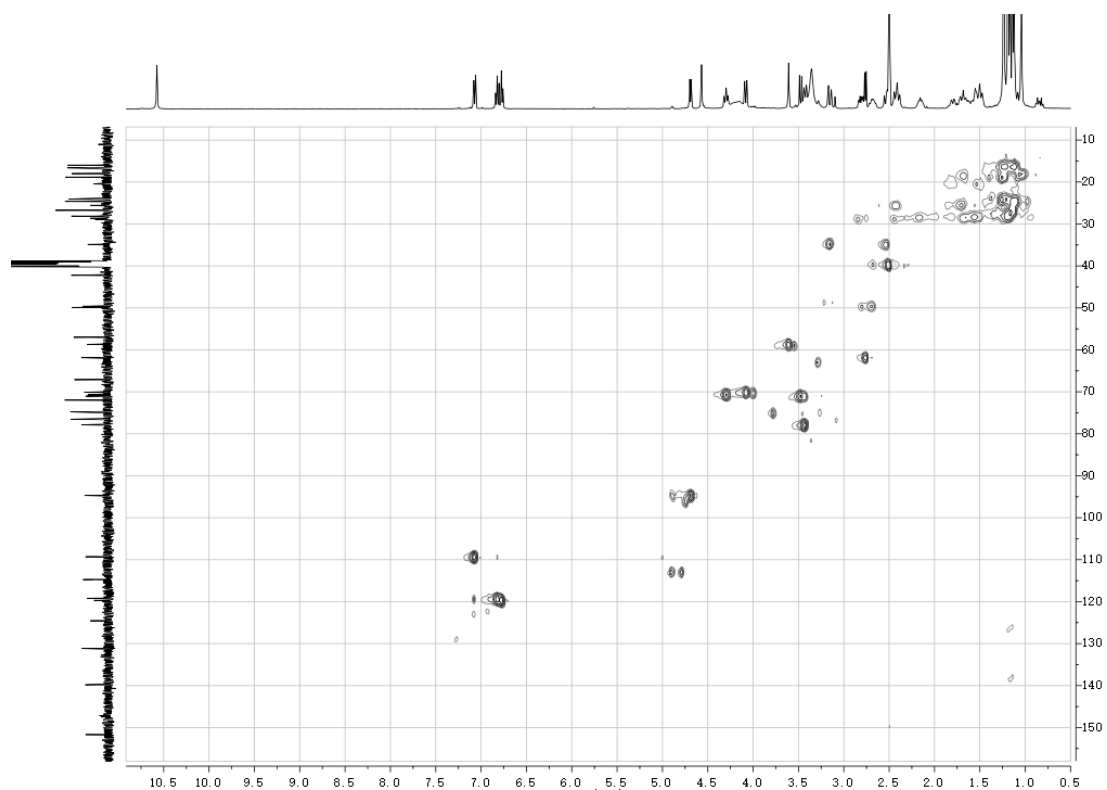


Figure S4: HSQC (100 MHz, DMSO- d_6) spectrum of **1** (Terpendole N)

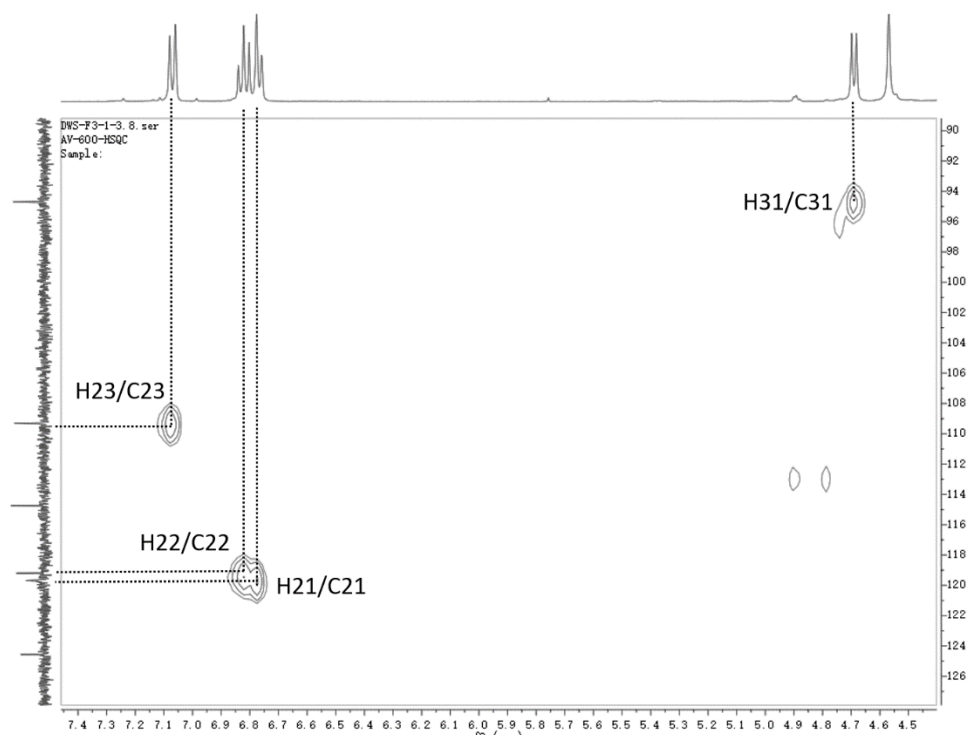


Figure S5: HSQC spectrum of **1** (Terpendole N) (From δ_C 90 ppm to δ_C 130 ppm)

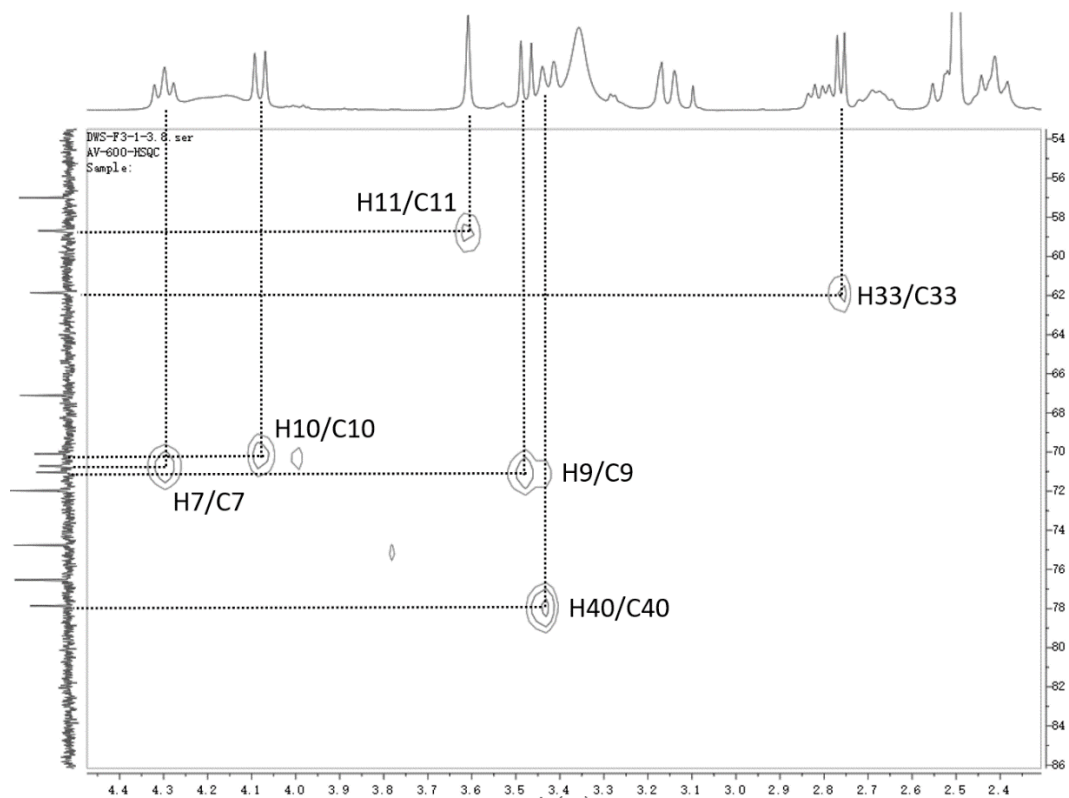


Figure S6: HSQC spectrum of **1** (Terpendole N) (From δ_C 54 ppm to δ_C 86 ppm)

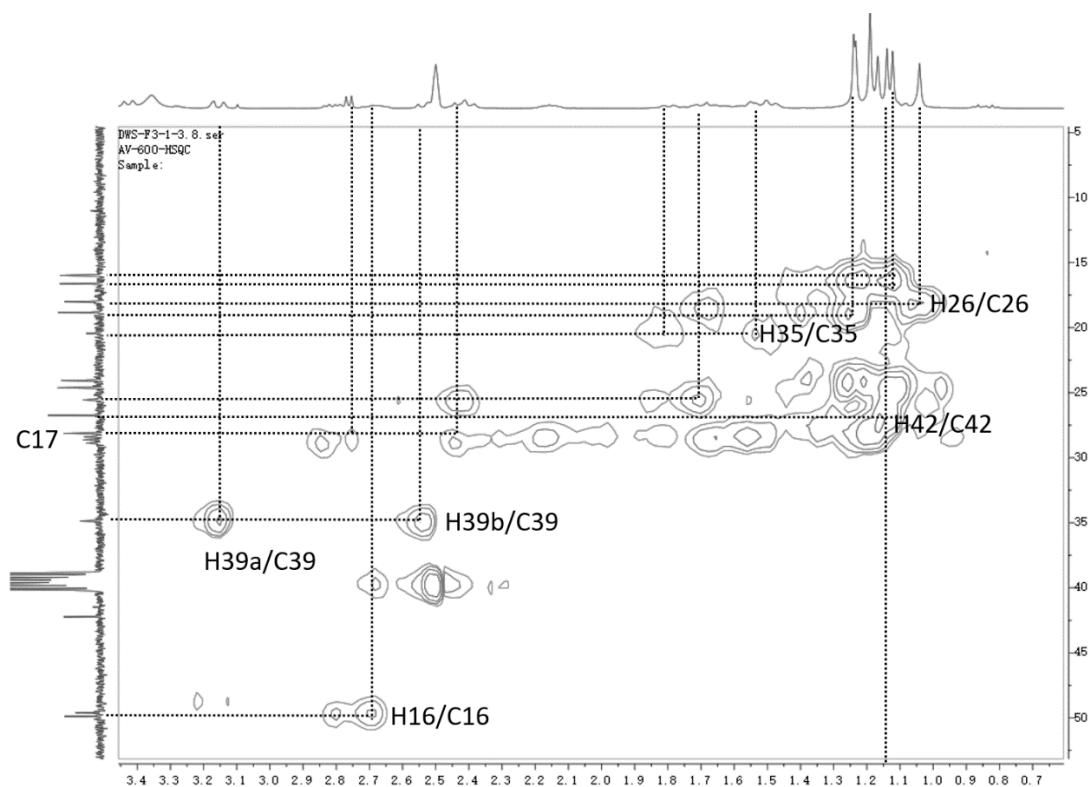


Figure S7: HSQC spectrum of **1** (Terpendole N) (From δ_C 5 ppm to δ_C 55 ppm)

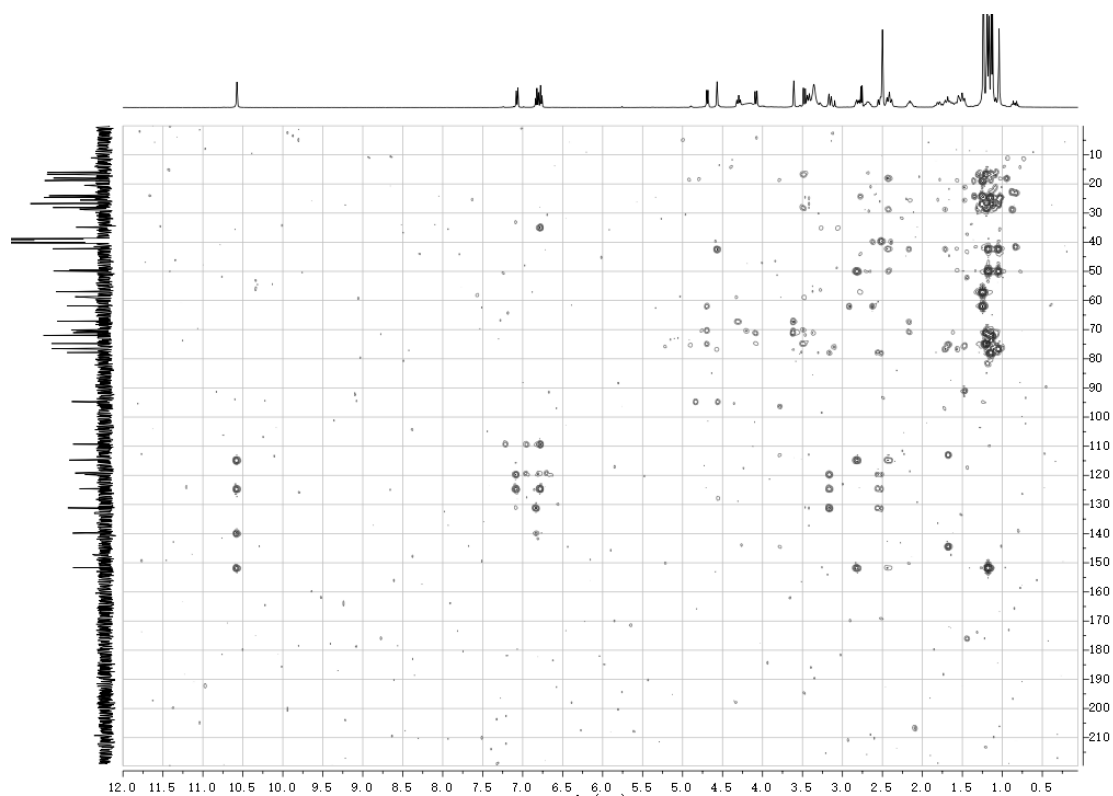


Figure S8: HMBC spectrum of **1** (Terpendole N)

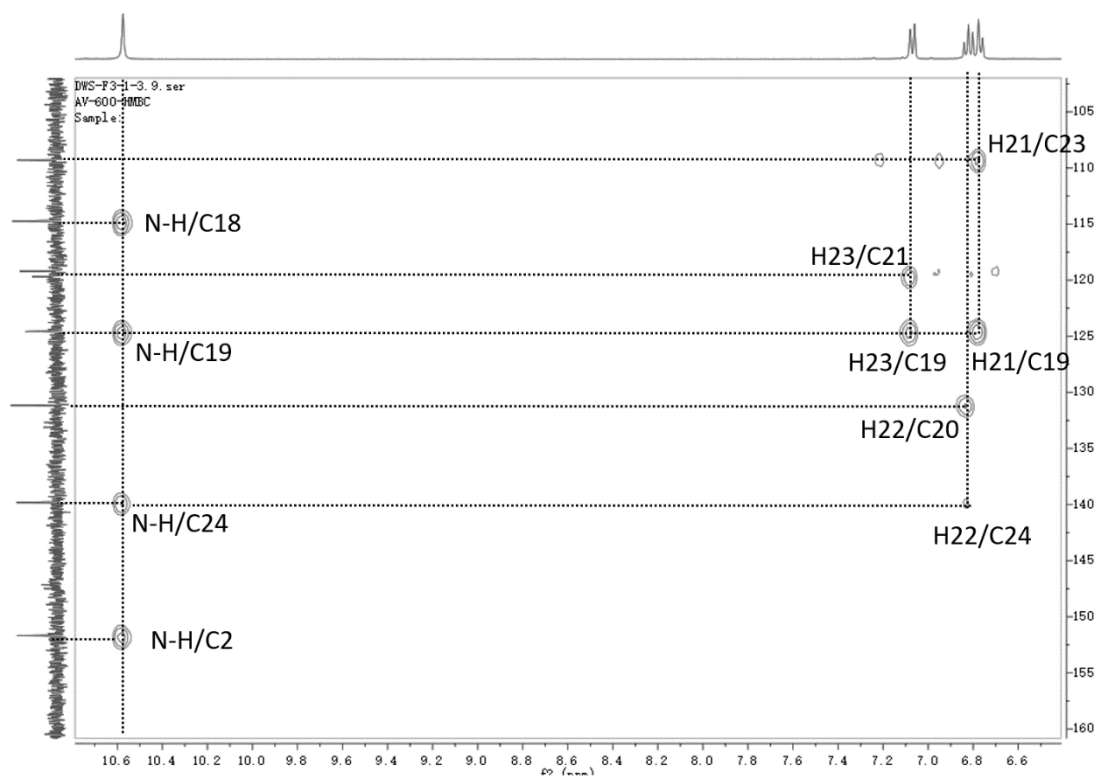


Figure S9: HMBC spectrum of **1** (Terpendole N) (From δ_{C} 100 ppm to δ_{C} 160 ppm)

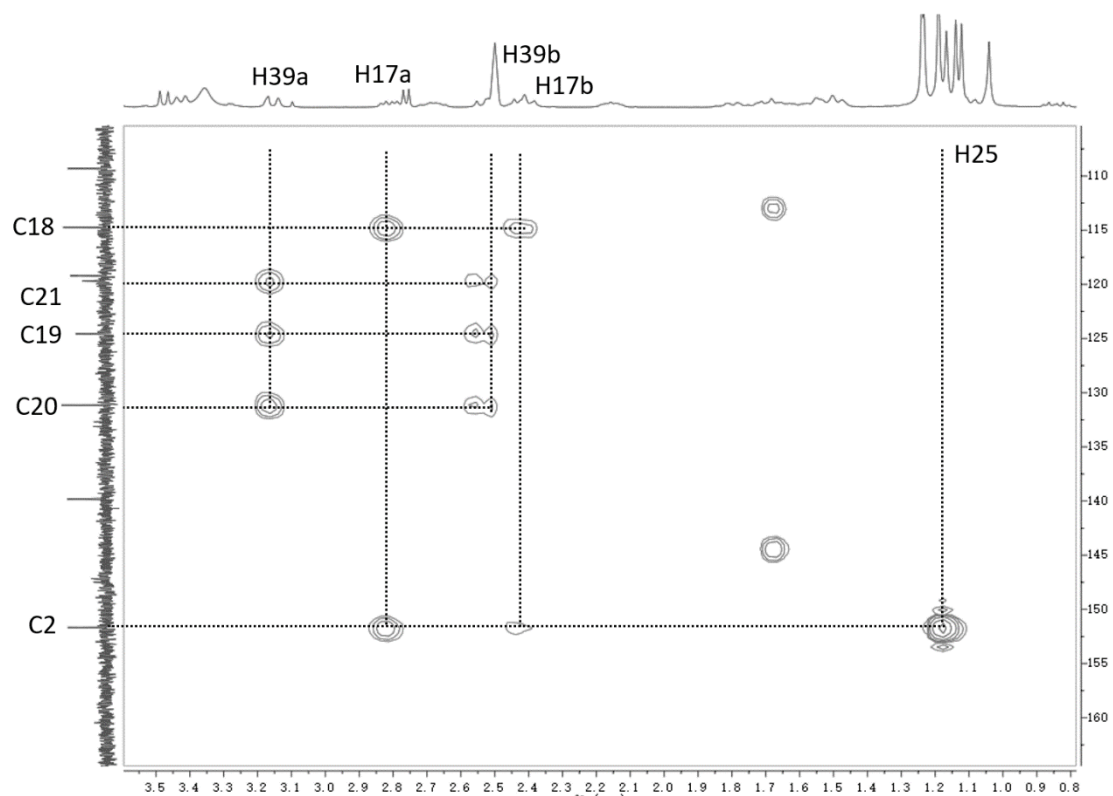


Figure S10: HMBC spectrum of **1** (Terpendole N) (From δ_{C} 100 ppm to δ_{C} 170 ppm)

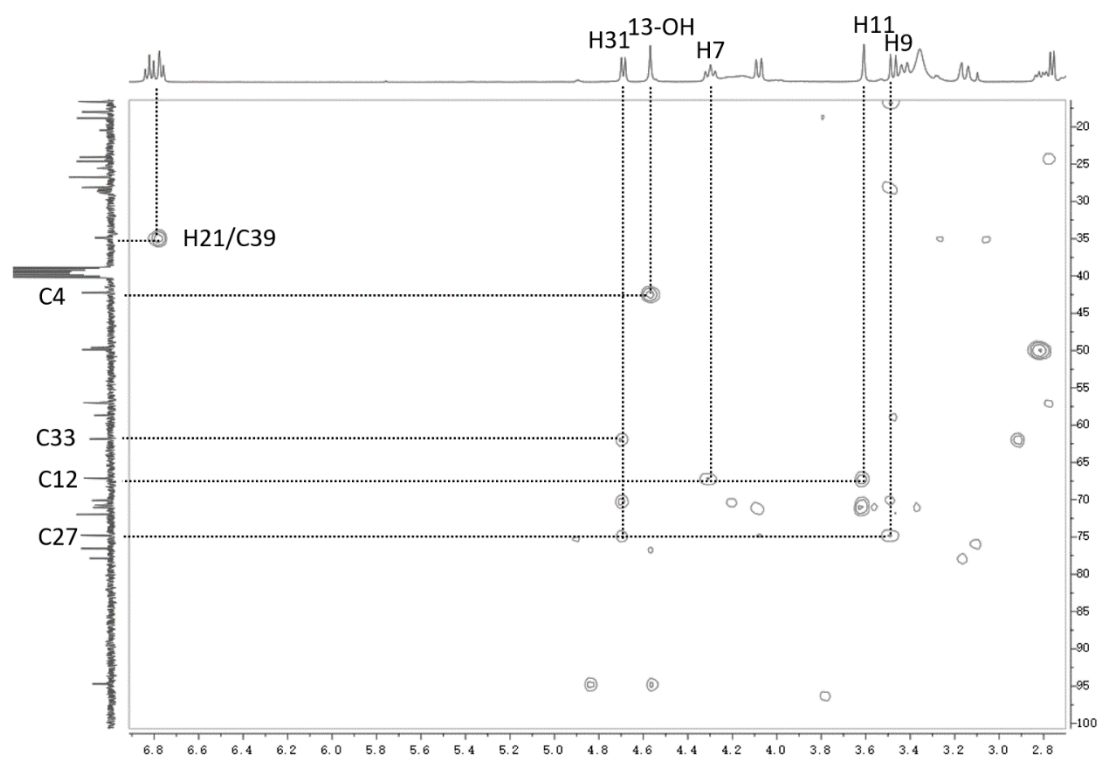


Figure S11: HMBC spectrum of **1** (Terpendole N) (From δ_{C} 20 ppm to δ_{C} 100 ppm)

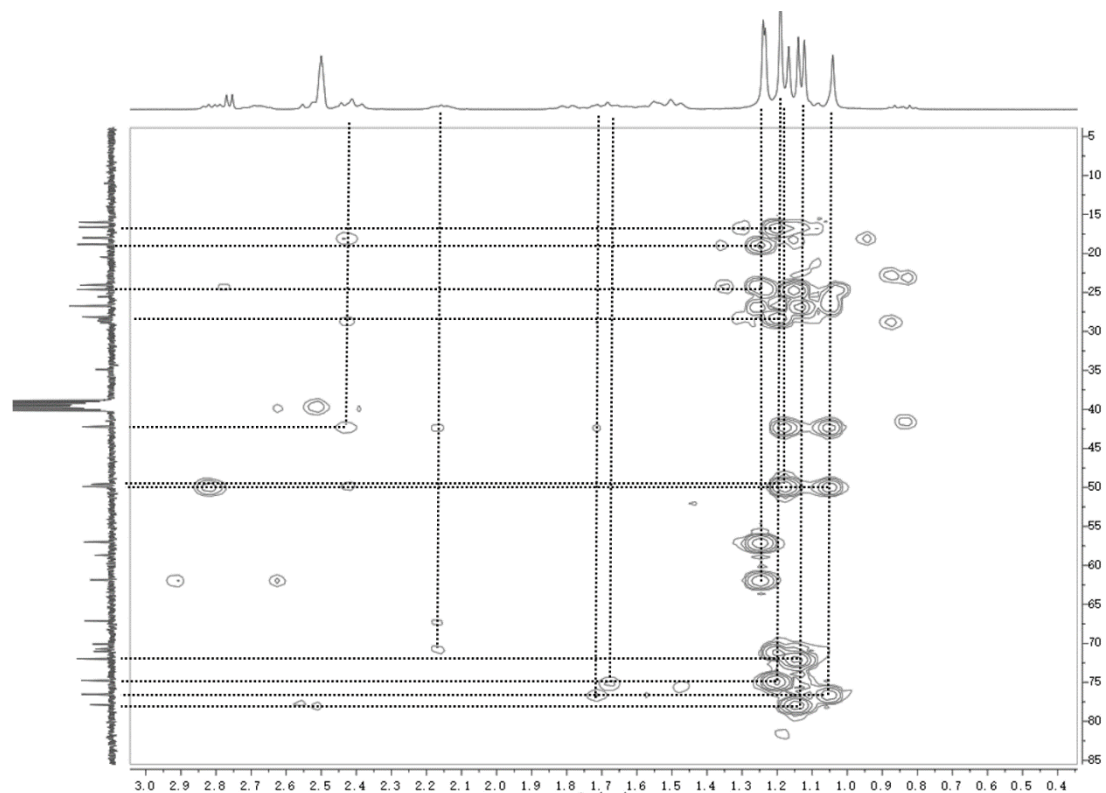


Figure S12: HMBC spectrum of **1** (Terpendole N) (From δ_{C} 5 ppm to δ_{C} 85 ppm)

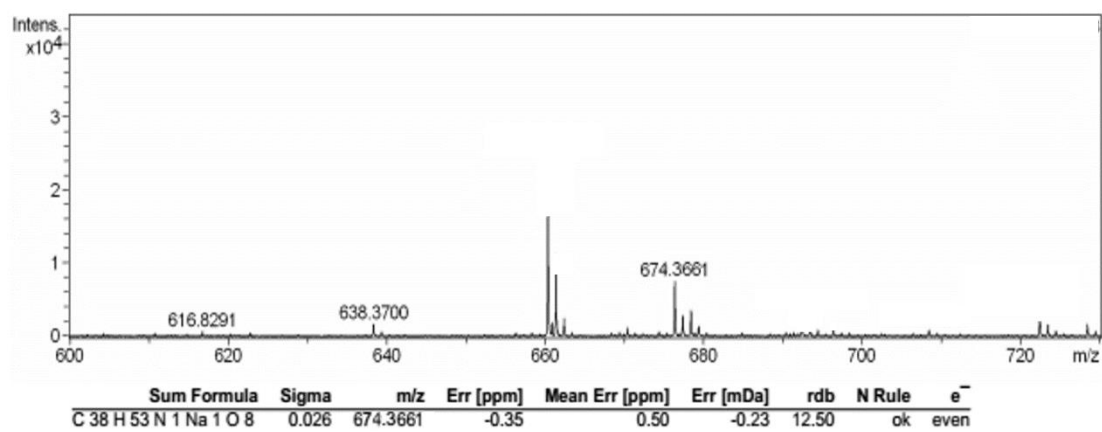


Figure S13: HR-ESI MS spectrum of **2** (Terpendole O)

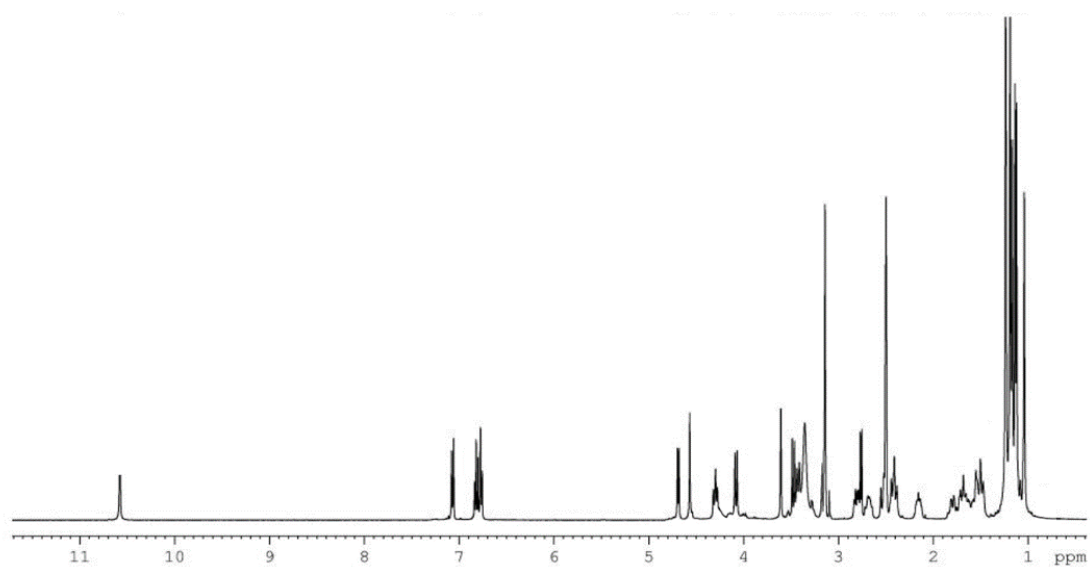


Figure S14: ^1H NMR (400 MHz, $\text{DMSO}-d_6$) spectrum of **2** (Terpendole O)

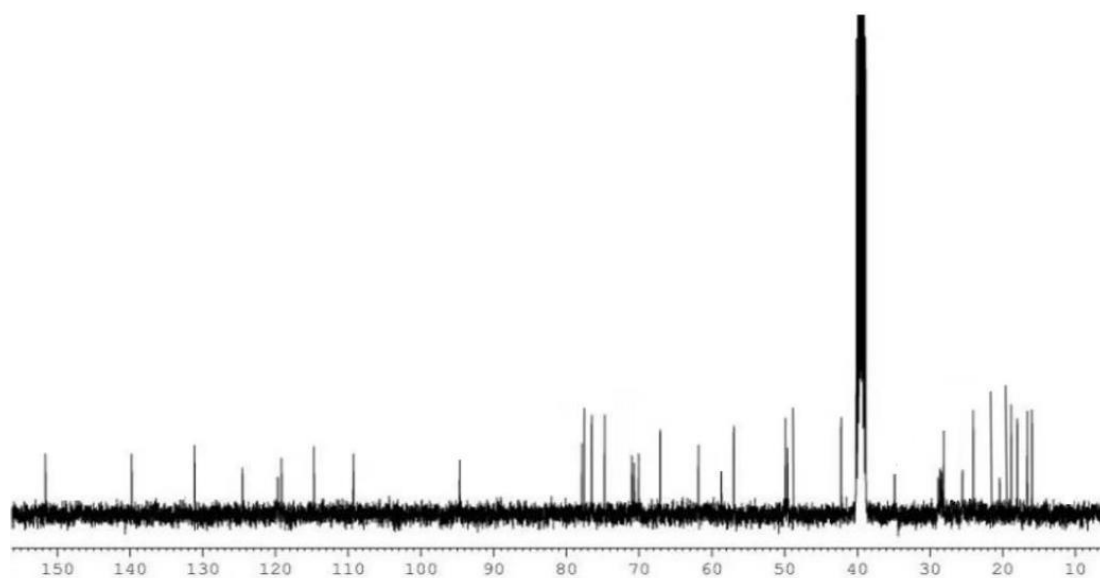


Figure S15: ^{13}C -NMR (100 MHz, $\text{DMSO-}d_6$) spectrum of **2** (Terpendole O)

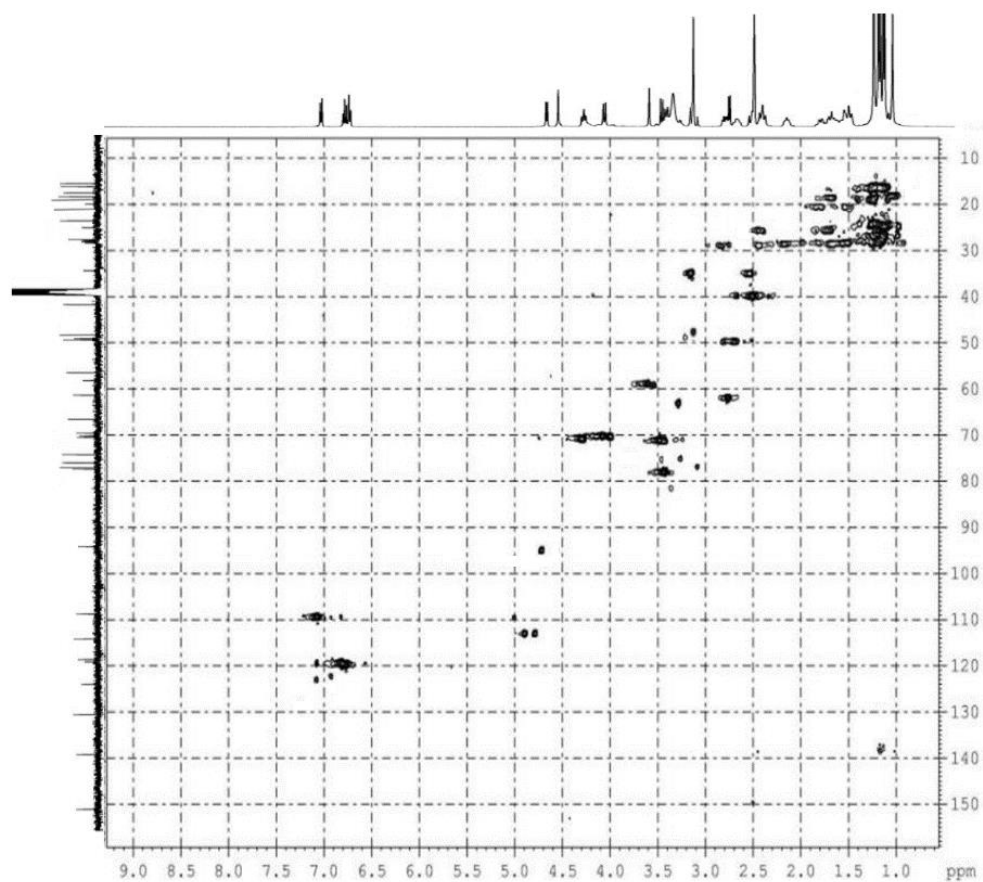


Figure S16: HSQC (100 MHz, DMSO- d_6) spectrum of **2** (Terpendole O)

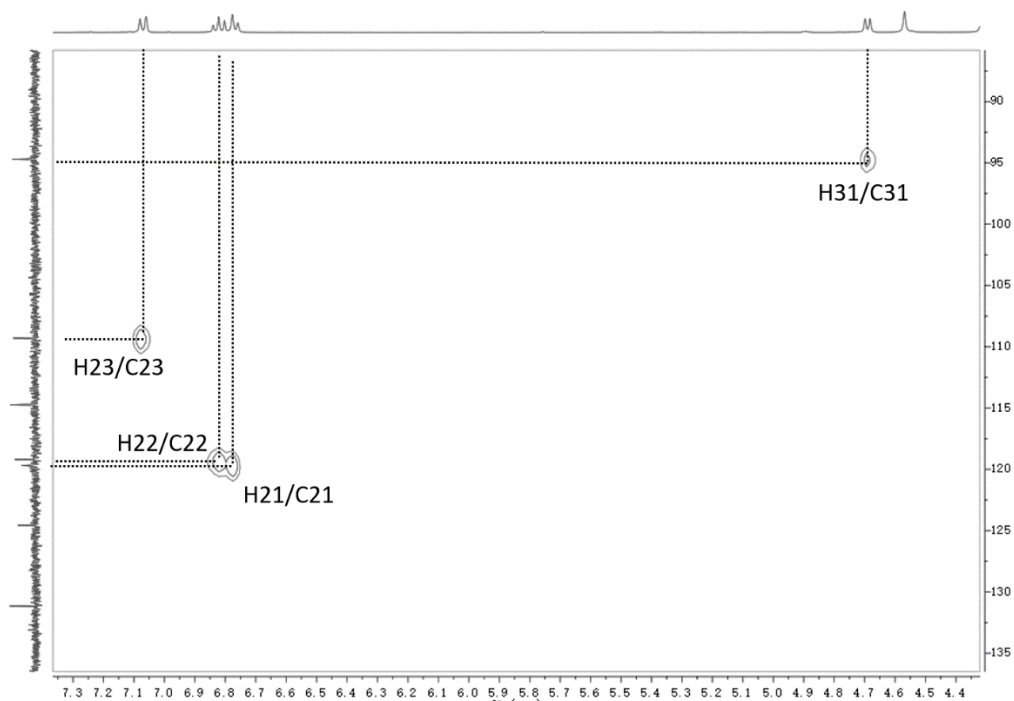


Figure S17: HSQC spectrum of **2** (Terpendole O) (From δ_{C} 80 ppm to δ_{C} 135 ppm)

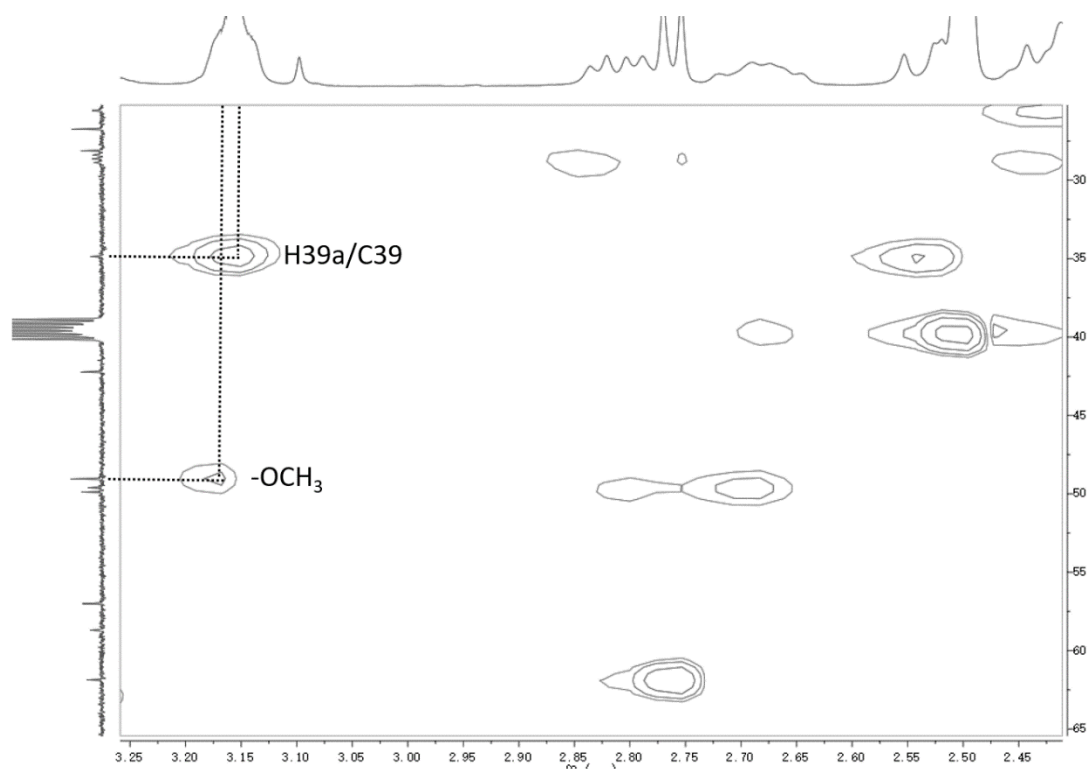


Figure S18: HSQC spectrum of **2** (Terpendole O) (From δ_{C} 20 ppm to δ_{C} 65 ppm)

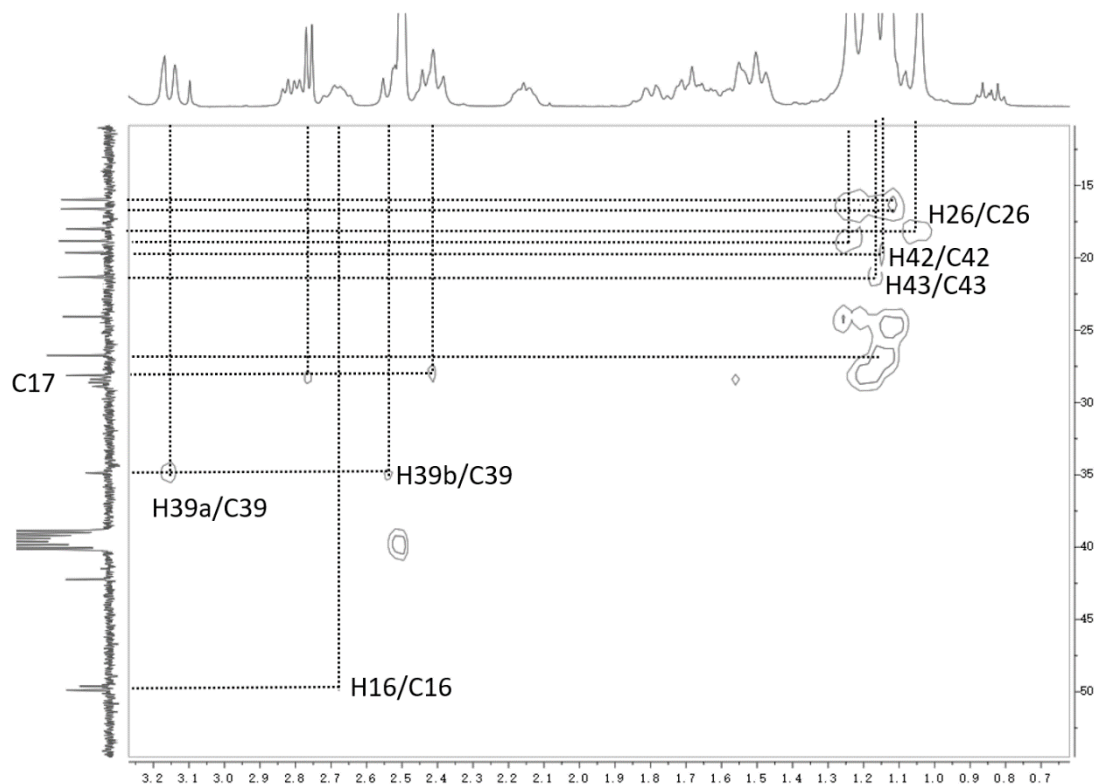


Figure S19: HSQC Spectrum of **2** (Terpendole O) (From δ_C 5 ppm to δ_C 55 ppm)

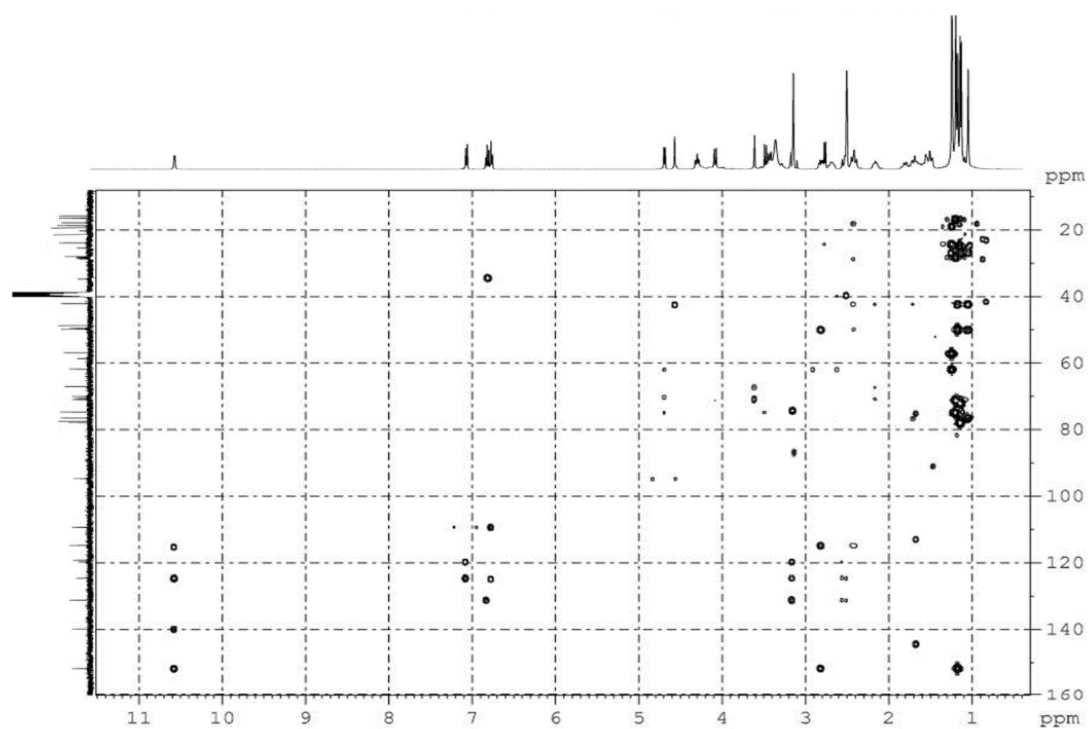


Figure S20: HMBC Spectrum of **2** (Terpendole O)

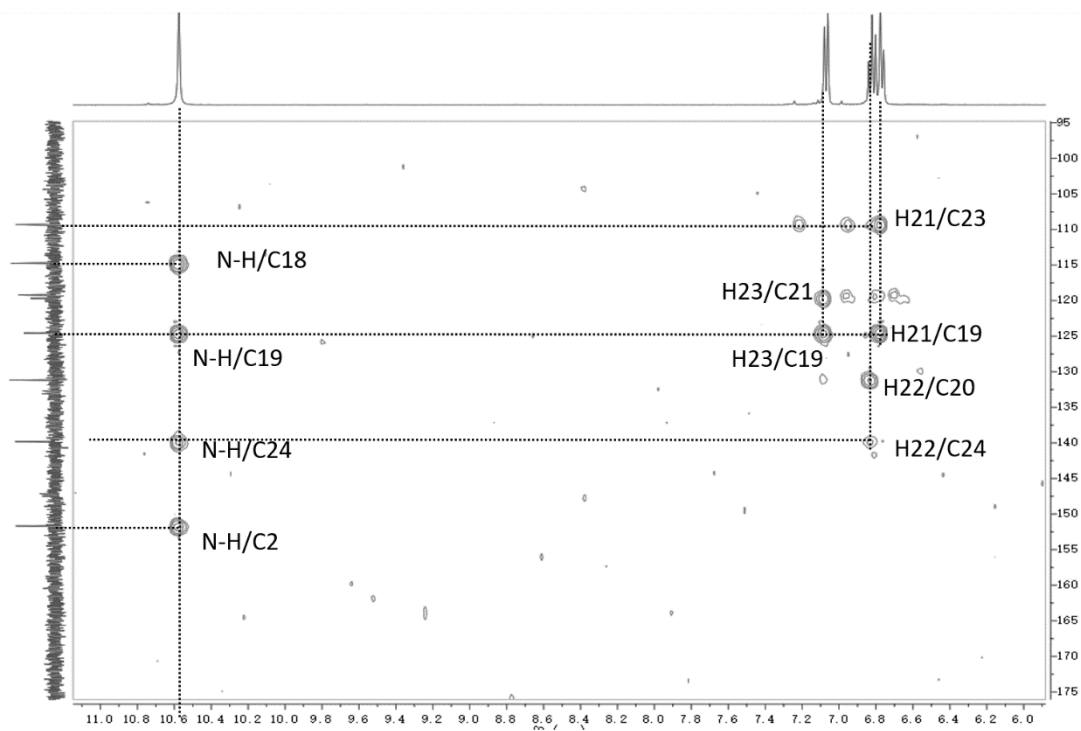


Figure S21: HMBC spectrum of **2** (Terpendole O) (From δ_C 95 ppm to δ_C 175 ppm)

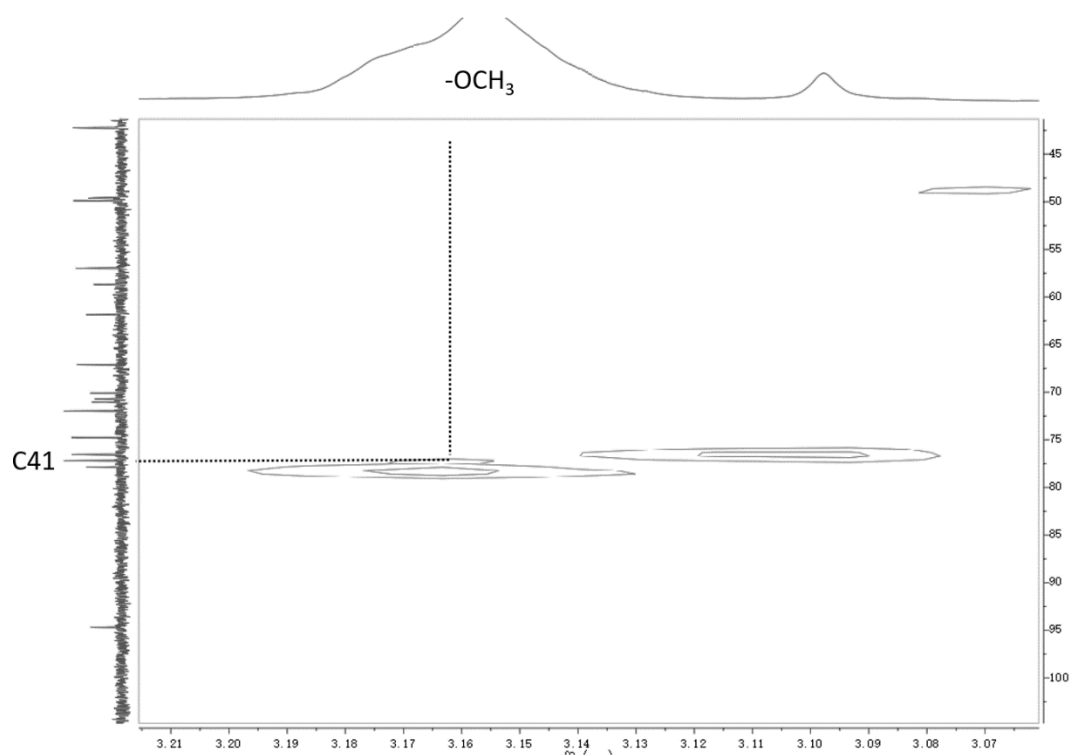


Figure S22: HMBC Spectrum of **2** (Terpendole O) (From δ_C 45 ppm to δ_C 100 ppm)

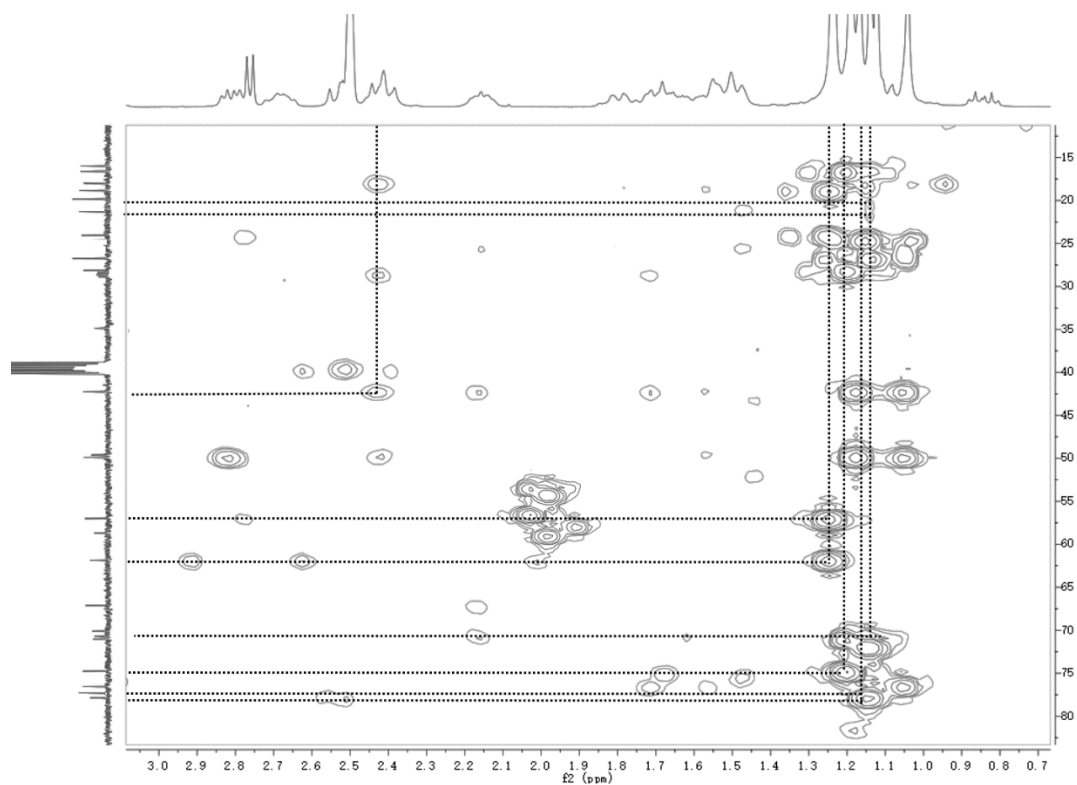
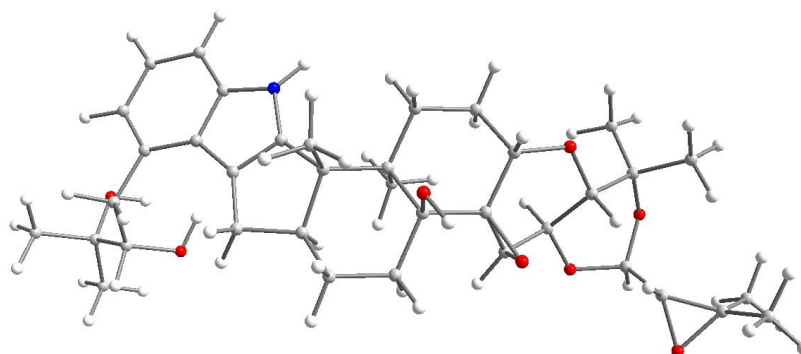


Figure S23: HMBC spectrum of **2** (Terpendole O) (From δ_{C} 10 ppm to δ_{C} 90 ppm)



X-ray Crystallographic Analysis of 1. The X-ray crystallographic data of **1** was collected on a Bruker APEX DUO diffractometer and equipped with an APEX II CCD using Mo K α radiation. **Crystal Data** for C₃₇H₅₁NO₈ ($M = 637.36$): triclinic, space group P-1 (no. 2), $a = 10.928(5)$ Å, $b = 11.701(5)$ Å, $c = 16.182(7)$ Å, $\alpha = 81.137(9)^\circ$, $\beta = 72.119(9)^\circ$, $\gamma = 80.132(9)^\circ$, $V = 1599.1(13)$ Å³, $Z = 2$, $T = 296.15$ K, $\mu(\text{Mo K}\alpha) = 0.097$ mm⁻¹, $D_{\text{calc}} = 1.299$ g/mm³, 16928 reflections measured ($2.88 \leq 2\theta \leq 50.7$), 5689 unique ($R_{\text{int}} = 0.0452$) which were used in all calculations. The final R_1 was 0.0603 ($>2\sigma(I)$) and wR_2 was 0.1513 (all data). 17101 reflections measured, 5799 unique ($R_{\text{int}} = 0.0461$) which were used in all calculations. The final $wR(F_2)$ was 0.1506 (all data). The crystal was kept at 296.15 K during data collection.

Figure S24: X-ray crystallographic analysis of **1**.

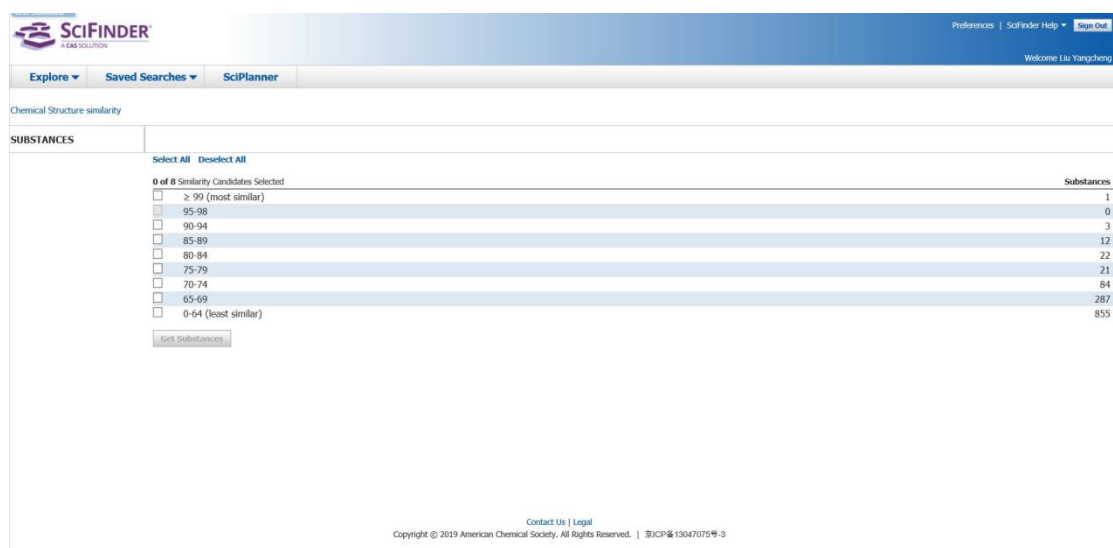


Figure S25: Scifinder search result of **1** (Similarity)

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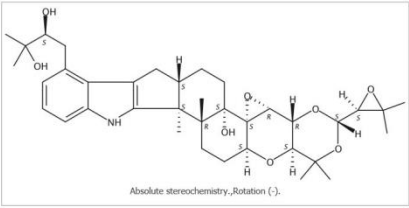
Link | Save | Print | Export

Chemical Structure similarity > substances (1) > 2337063-57-5

SUBSTANCE DETAIL | Get References | Send to SciPlanner

Return

CAS Registry Number 2337063-57-5



$C_{27}H_{41}NO_8$
INDEX NAME NOT YET ASSIGNED

Molecular Weight
637.80

Density (Predicted)
Value: 1.3540.1 g/cm³ | Condition: Temp: 20 °C Press: 760 Torr

pKa (Predicted)
Value: 14.6040.70 | Condition: Most Acidic Temp: 25 °C

Other Names
Tolyposcadin I

Absolute stereochemistry, Rotation (-).

Expand All | Collapse All

▶ PREDICTED PROPERTIES
▶ PREDICTED SPECTRA
▶ BIOACTIVITY INDICATORS

Figure S26: Scifinder search result of **1** (Substance Detail)

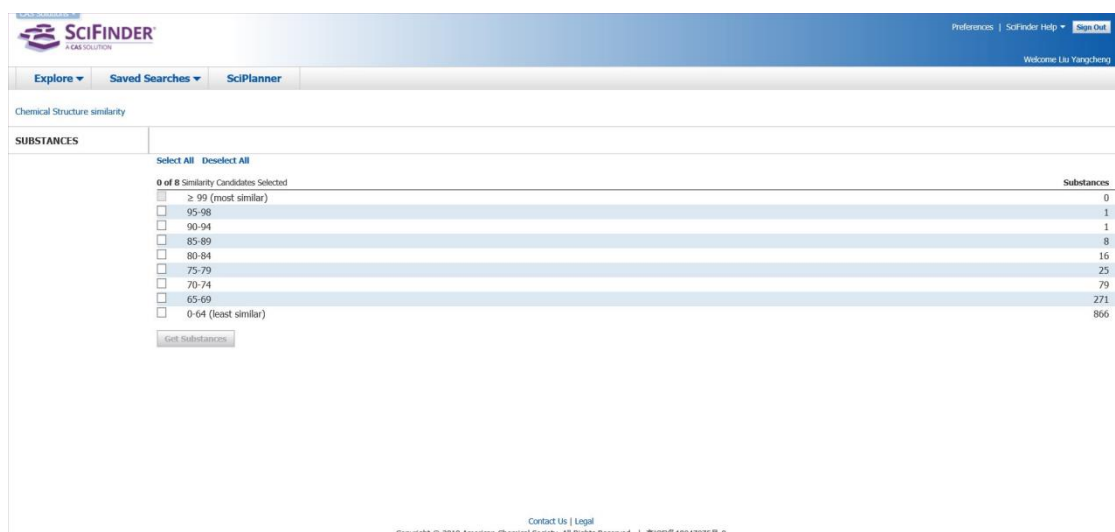


Figure S27: Scifinder search result of 2 (Similarity)

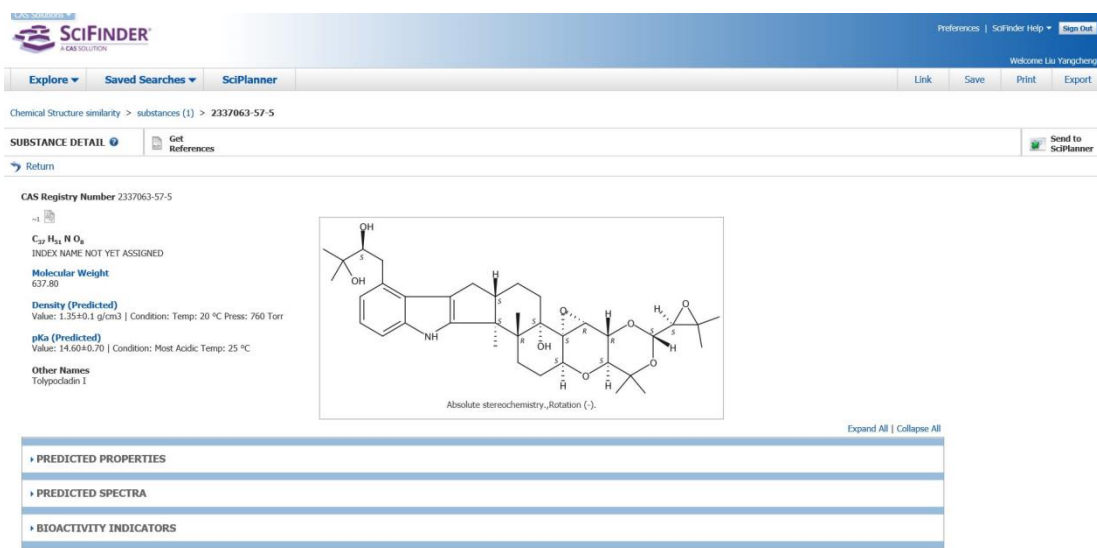


Figure S28: Scifinder search result of **2** (Substance Detail)