

Supplementary Data

N-(Isobutyl)-3,4-methylenedioxy Cinnamoyl Amide

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Figure S1. High resolution mass spectrometry ion chromatogram for FD fraction of *Zanthoxylum zanthoxyloides*.

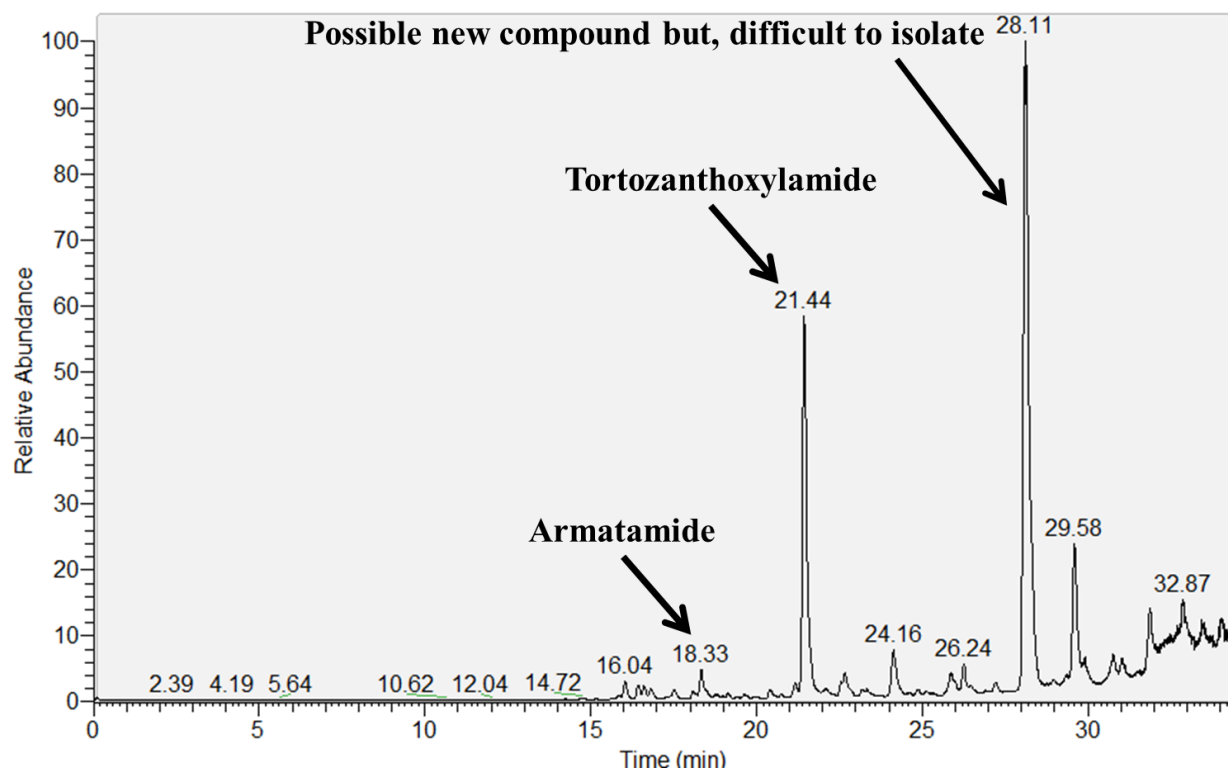


Figure S2. Mass spectrometry fragmentation ions for Armatamide.

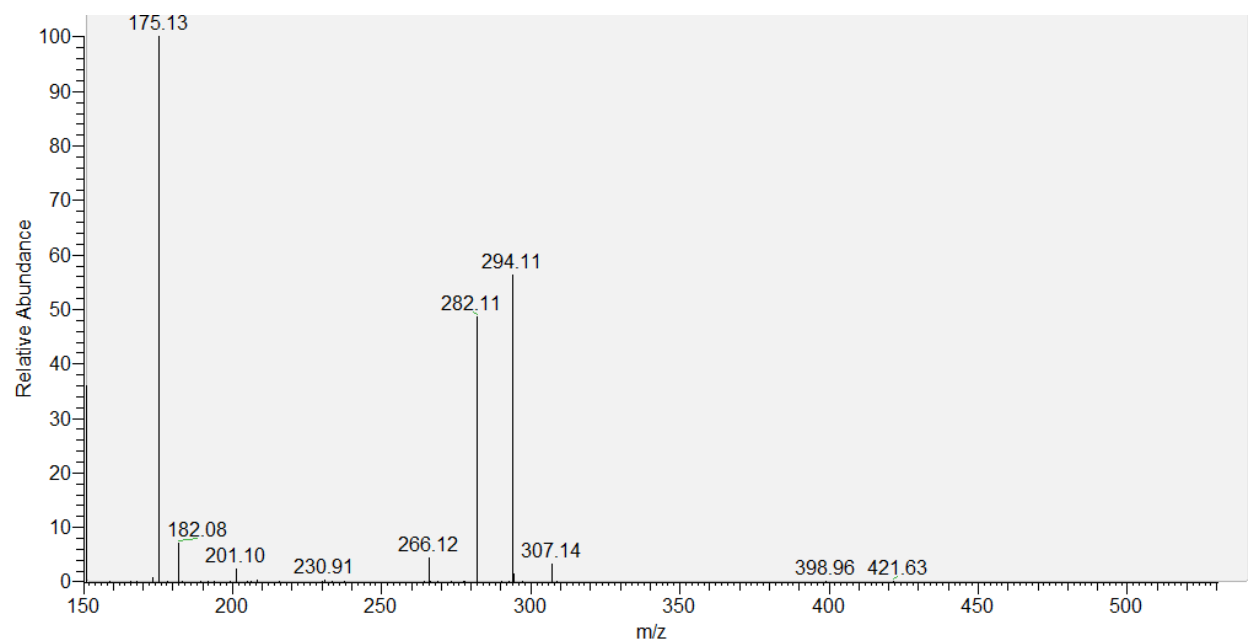


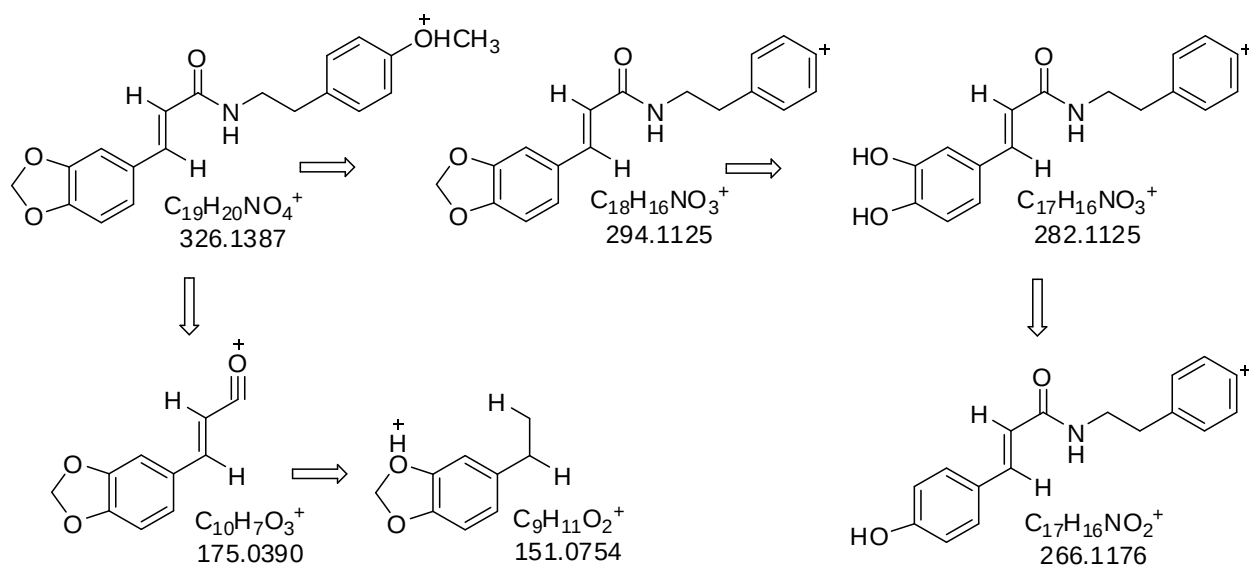
Figure S3. Proposed mass spectrometry fragmentation structures for Armatamide.

Figure S4. High resolution mass spectrometry ion chromatogram for tortozanthoxylamide (**1**).

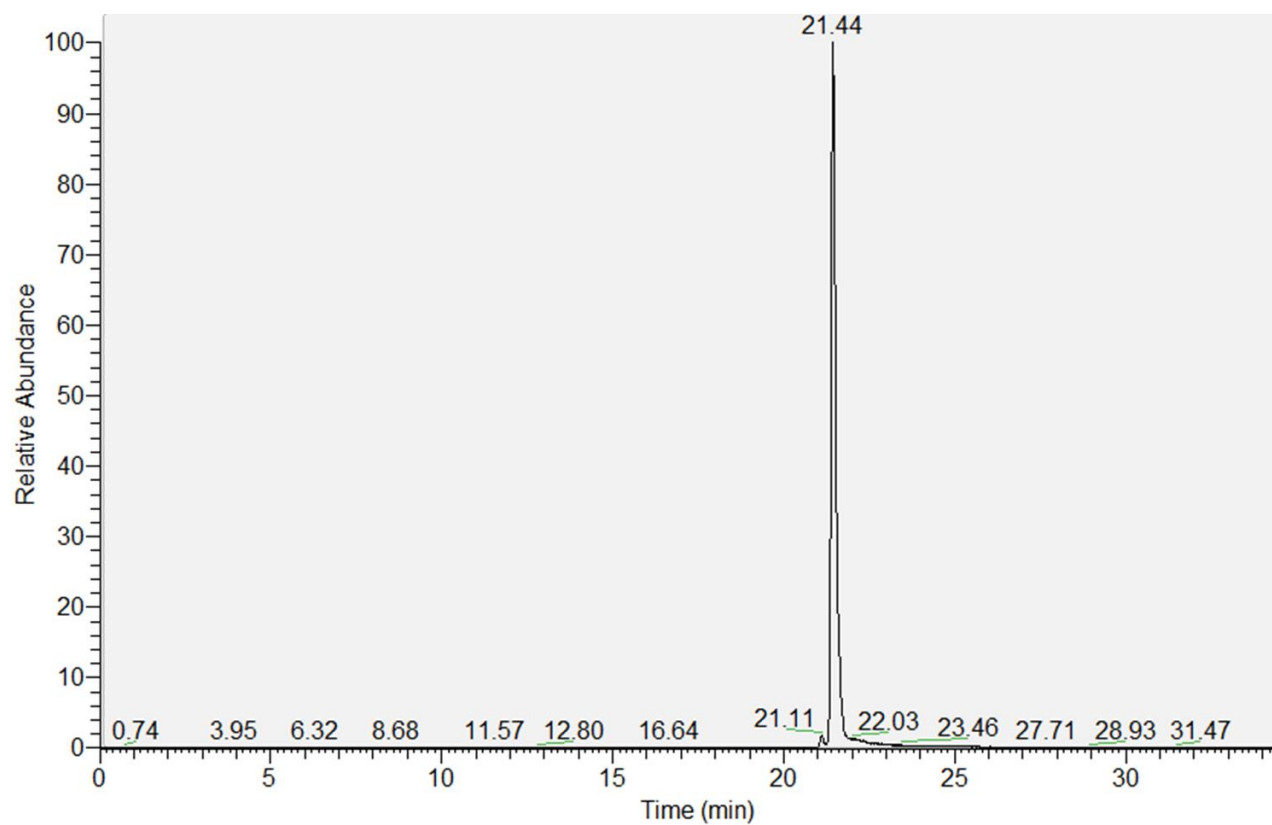


Figure S5. Mass spectrometry fragmentation ions for tortozanthoxylamide.

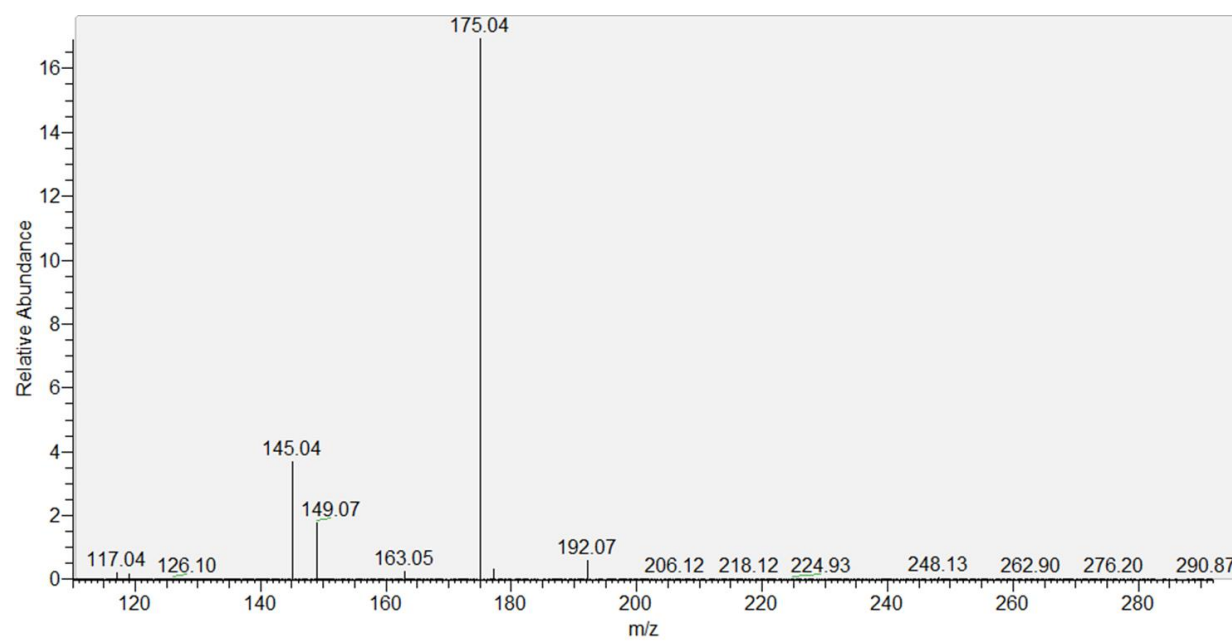


Figure S6. Proposed mass spectrometry fragmentation structures for tortozanthoxylamide.

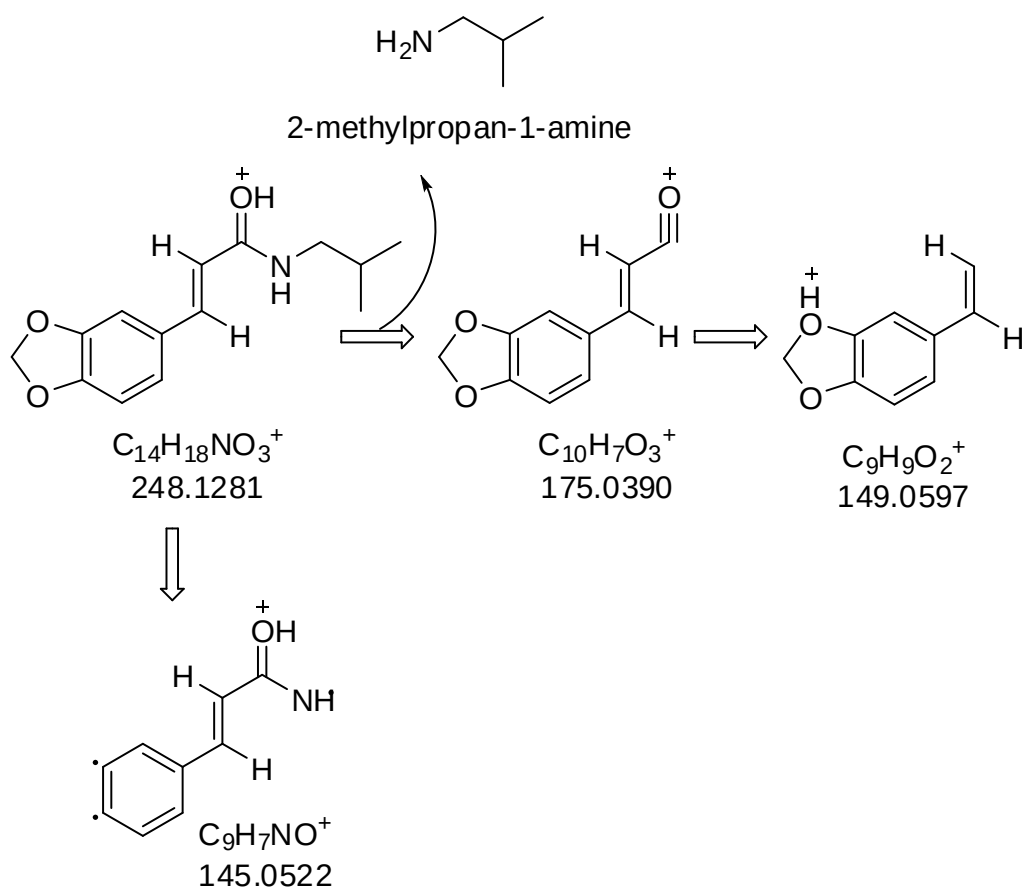


Figure S7. Mass spectrometry fragmentation ions for lanyuamide I-III analogue.

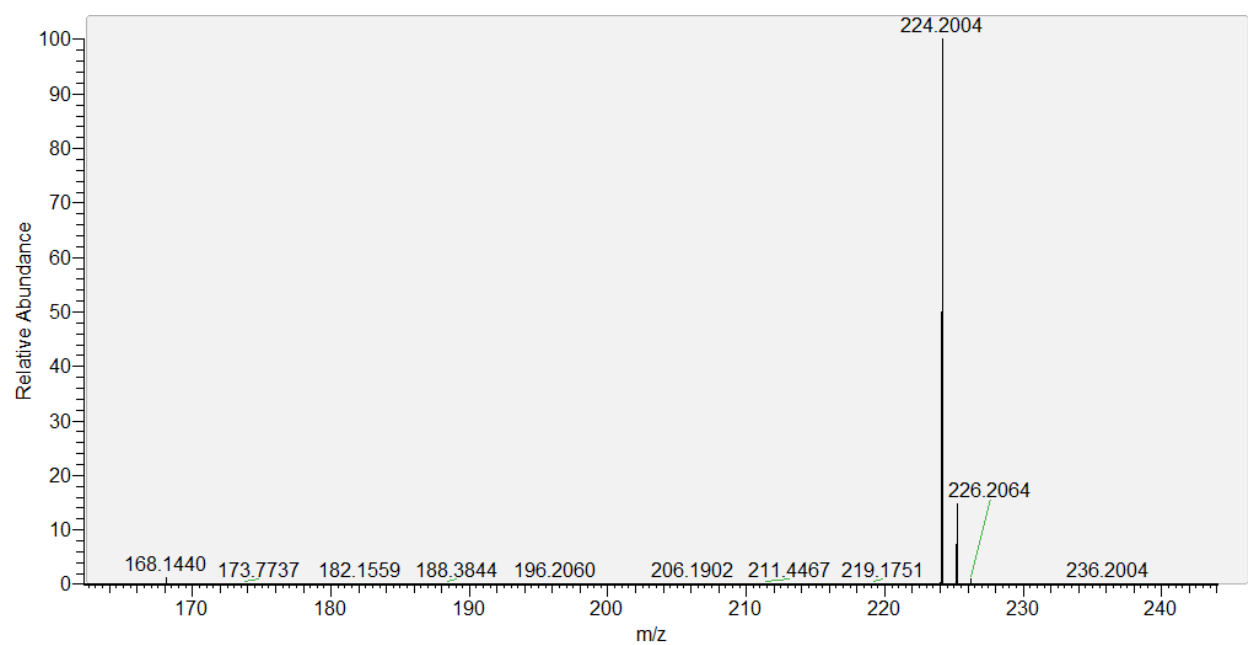


Figure S8. Proposed mass spectrometry fragmentation structures for lanyuamide I-III analogue.

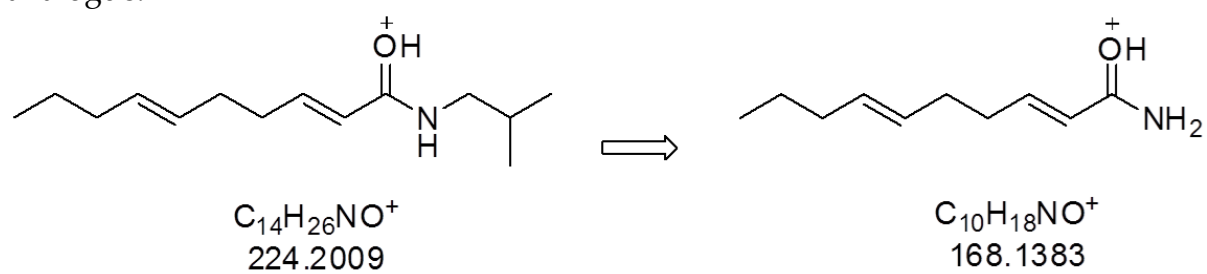


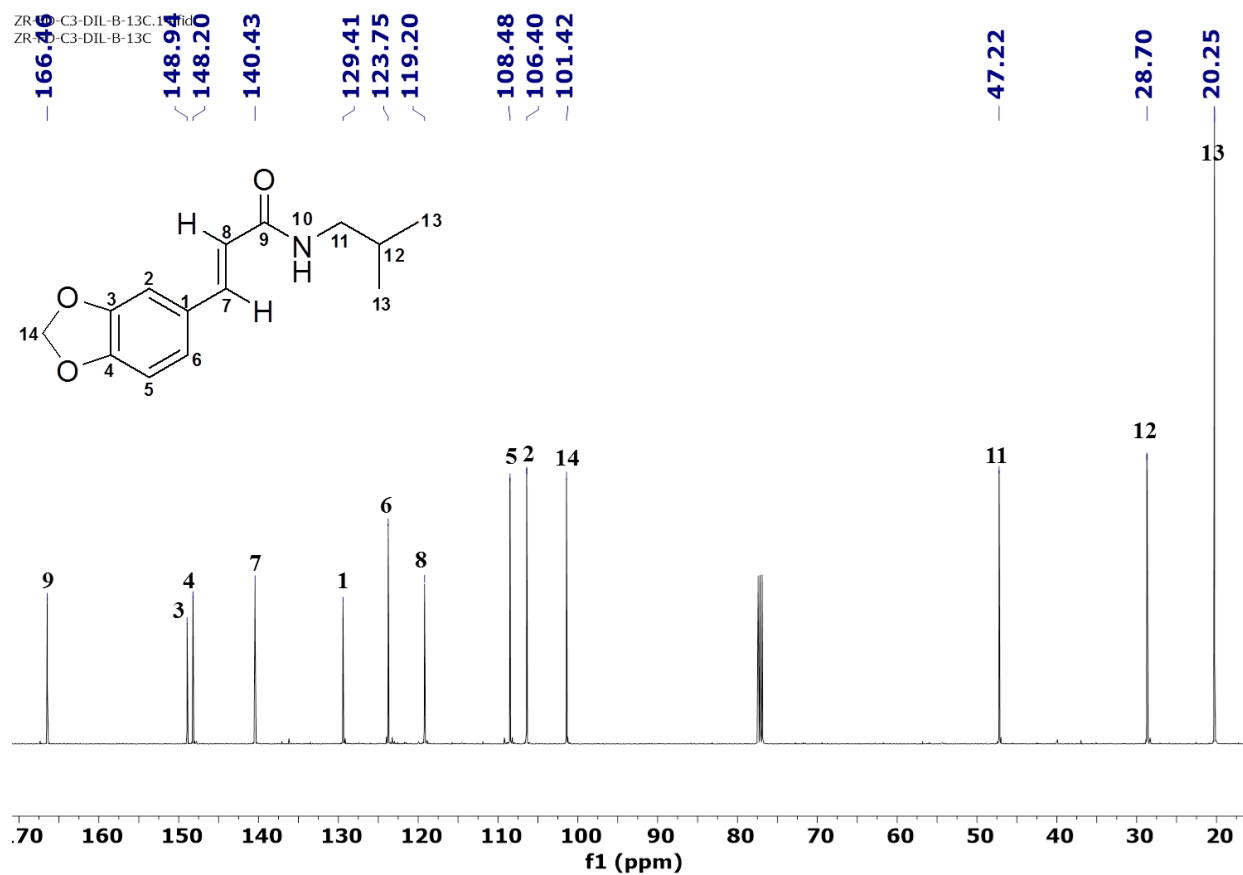
Figure S9. ^{13}C -NMR spectrum of tortozanthoxylamide (1) in CDCl_3 .

Figure S10. ^1H -NMR spectrum of tortozanthoxylamide (**1**) in CDCl_3 .

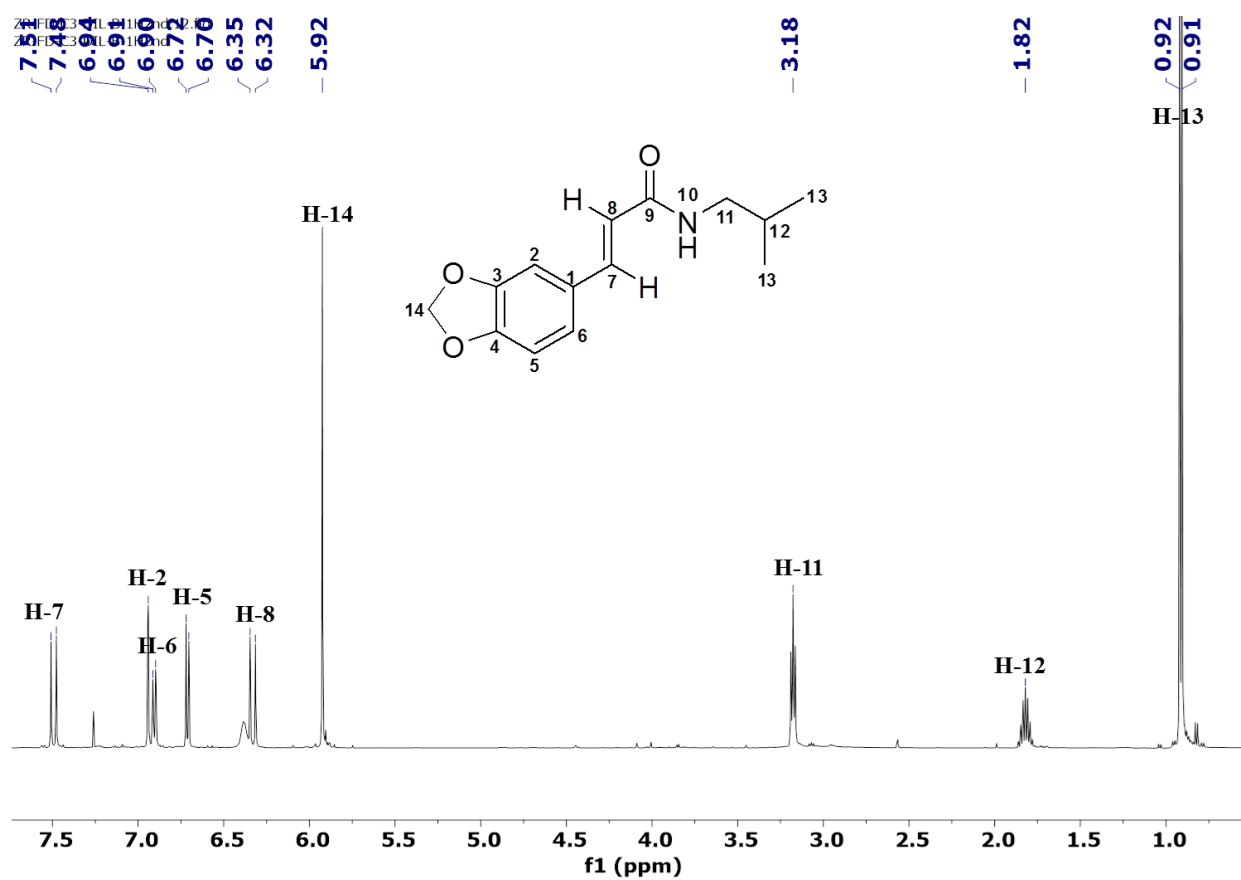


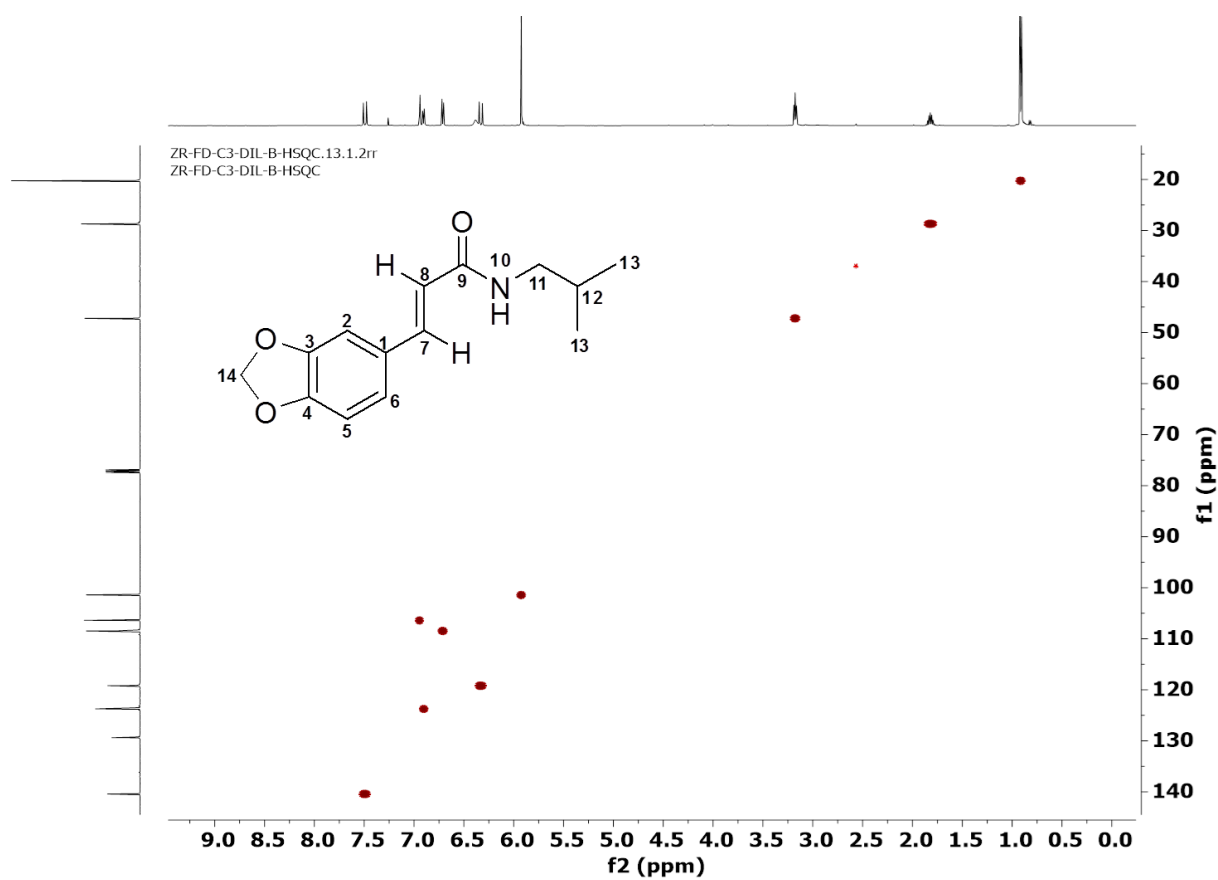
Figure S11. HSQC spectrum of tortozanthoxylamide (**1**) in CDCl₃.

Figure S12. COSY spectrum of tortozanthoxylamide (**1**) in CDCl₃.

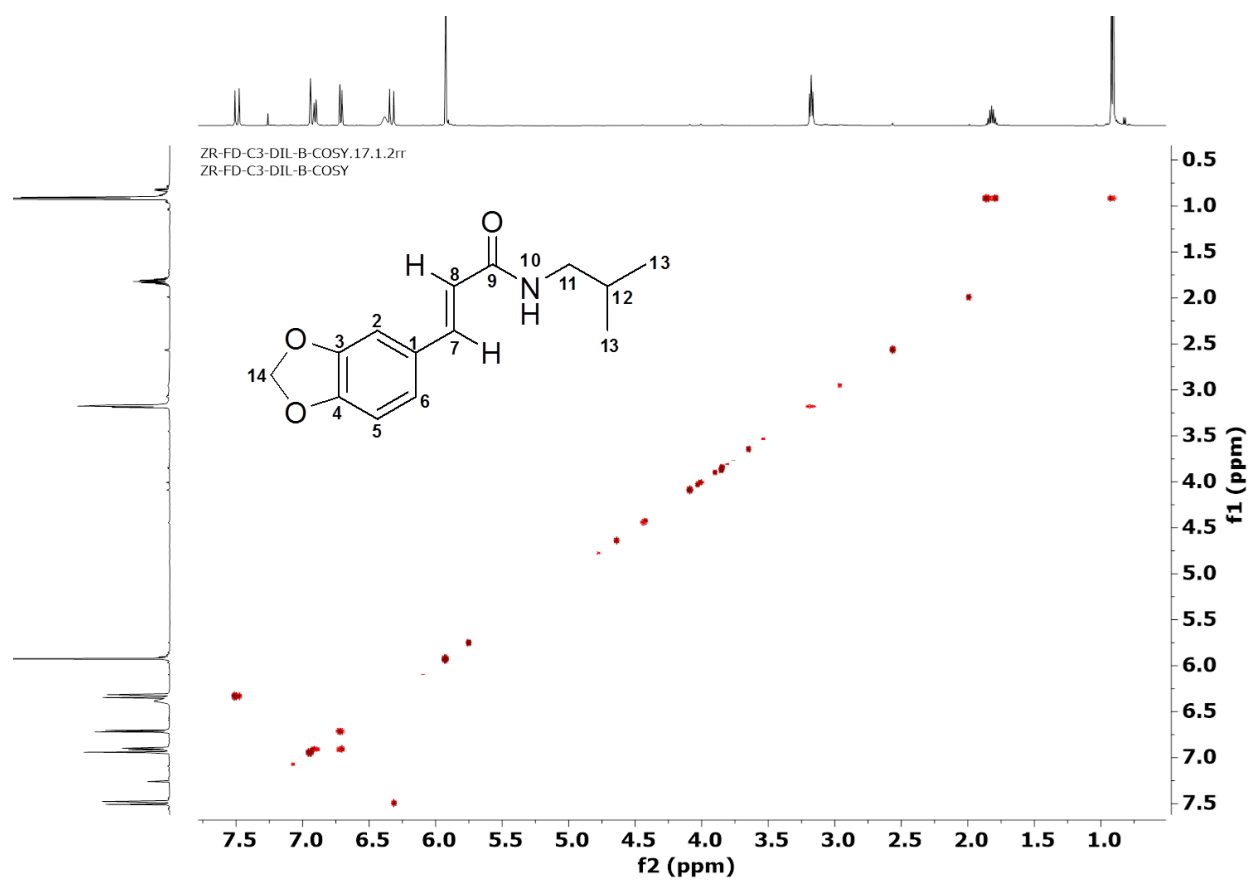


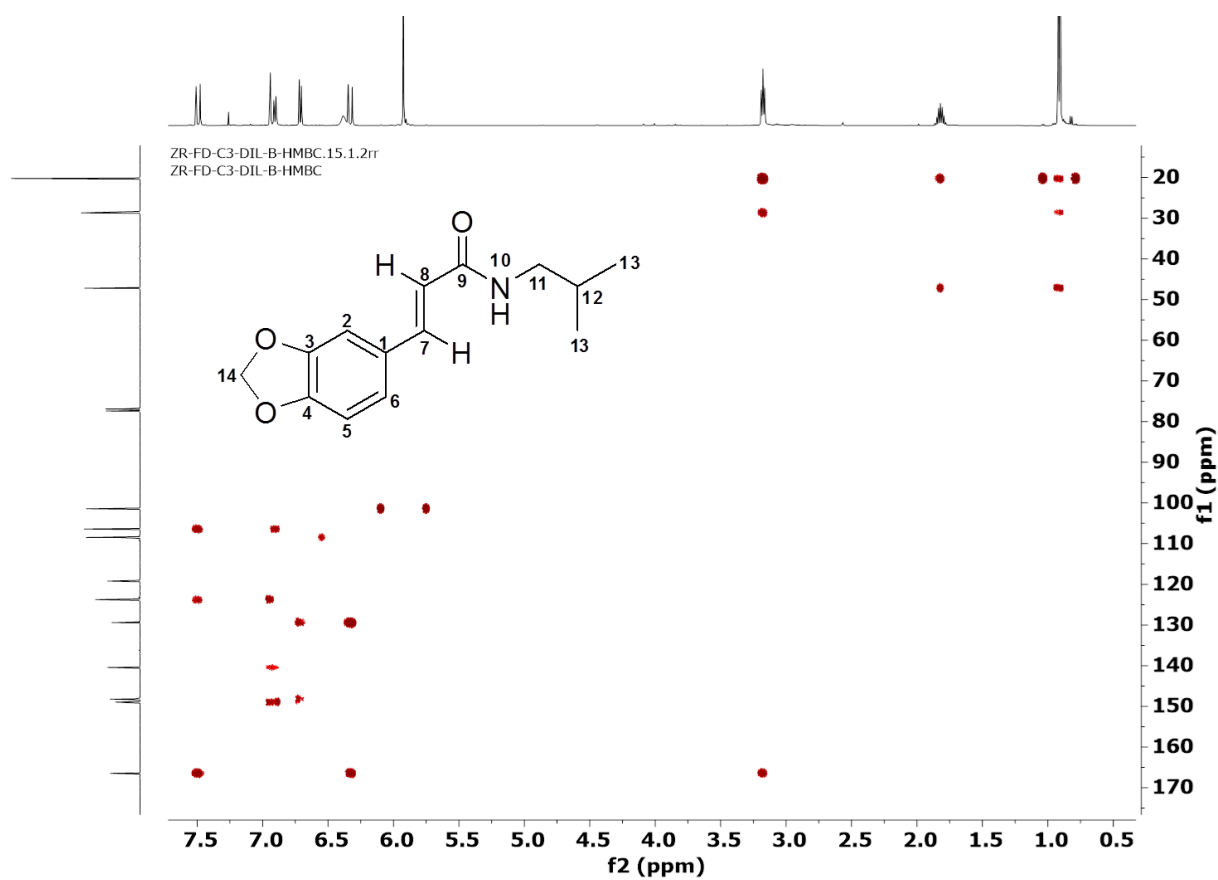
Figure S13. HMBC spectrum of tortozanthoxylamide (1) in CDCl₃.

Figure S14. TOCSY spectrum of tortozanthoxylamide (**1**) in CDCl₃.

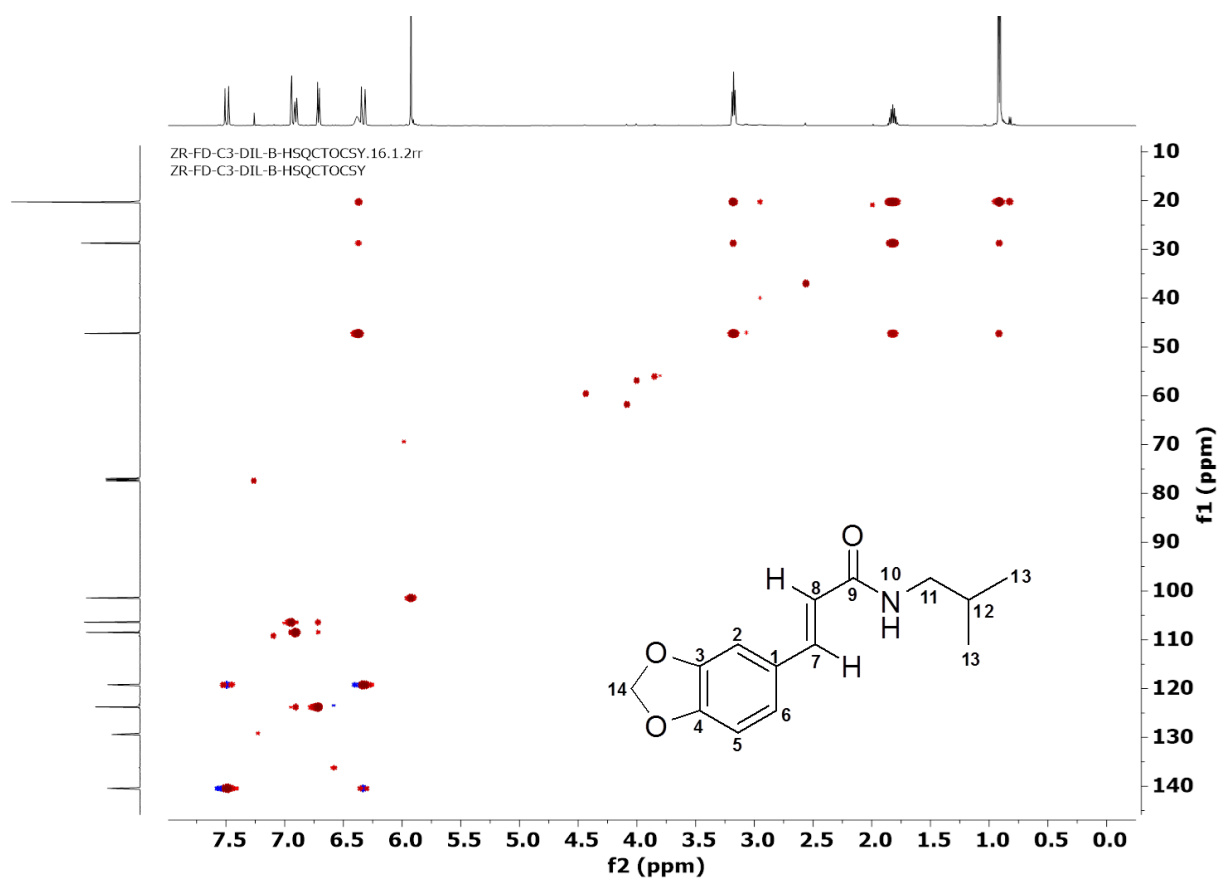


Figure S15. T-ROESY spectrum of tortozanthoxylamide (**1**) in CDCl₃.

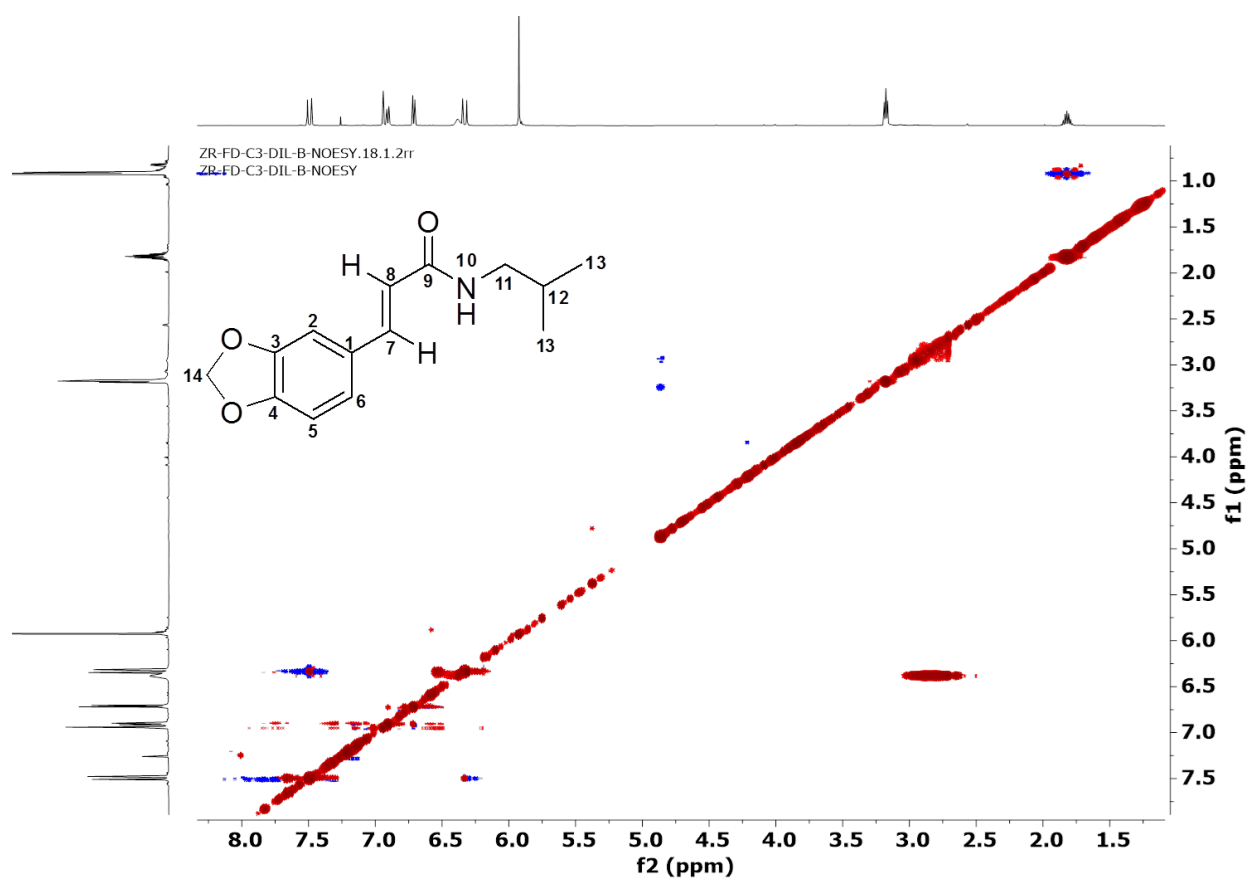


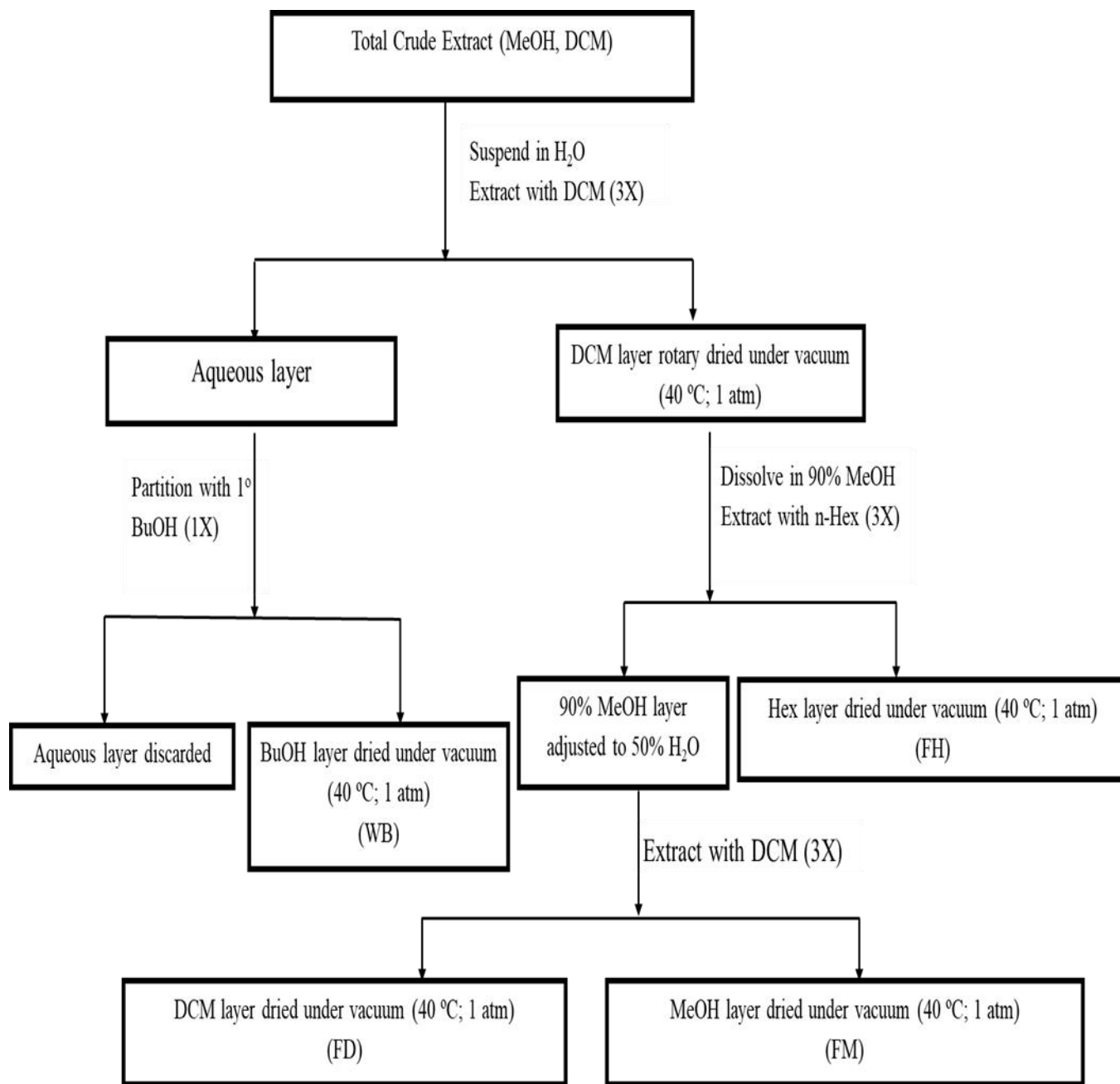
Figure S16. Flow chart of the modified Kupchan solvent partitioning process.

Table S1. Mean percentage cell count at cell cycle phases of *T. brucei* subsp. *brucei*.

COMPOUNDS	PERCENTAGE CELL COUNT (MEAN $\pm$ SEM)		
	G0/G1 Phase	S Phase	G2/M Phase
Negative control	62.2 $\pm$ 0.7	13.1 $\pm$ 0.2	24 $\pm$ 1
Tortozanthoxylamide	56.2 $\pm$ 0.9	19.4 $\pm$ 0.9	26.8 $\pm$ 0.4

Table S2. Effects of FD fractions on cell viability of *T. brucei* subsp. *brucei*.

FD, Subsequent fractions, and Standard	MEAN IC ₅₀ $\pm$ SE (μ g/ml)
FD	7.9 $\pm$ 0.1
C1	11.2 $\pm$ 0.7
C2	13.4 $\pm$ 0.5
C3	9 $\pm$ 1
C4	2.0 $\pm$ 0.2
C5	6 $\pm$ 1
C6	10 $\pm$ 1
C7	17 $\pm$ 3
C8	8.9 $\pm$ 0.7
C9	19 $\pm$ 2
C10	28 $\pm$ 1
C11	54 $\pm$ 8
DA	0.53 $\pm$ 0.04

Mean IC₅₀ and standard errors (SE) were calculated from three different experiments. DA=Diminazene aceturate was used as a positive control.