

SUPPLEMENT MATERIAL

Bioassay-guided isolation and HPLC quantification of antiproliferative metabolites from *Stahlianthus thorelii*

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¹H NMR and ¹³C-NMR spectrums, and HREIMS of *Stahlianthus thorelli*

1. ¹H NMR and ¹³C-NMR spectrums, and HREIMS of (E)-1-(3',5'-dihydroxy-4'-(2''-hydroxybenzyl) -2'- methoxyphenyl) -3- (2,4-dimethoxyphenyl) prop-2-en-1-one (1)

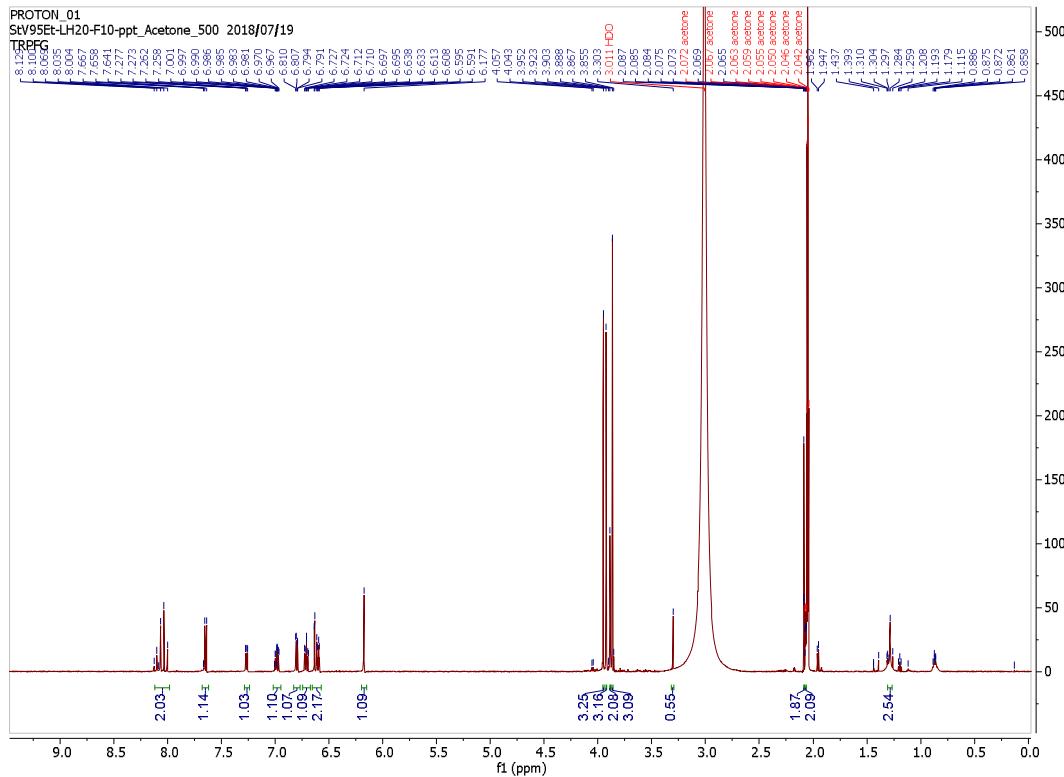


Figure S1. ¹H NMR spectrum of ((E)-1-(3',5'-dihydroxy-4'-(2''- hydroxybenzyl) -2'- methoxyphenyl) -3- (2,4-dimethoxyphenyl) prop-2-en-1-one) (**1**) measured in 500 MHz in acetone

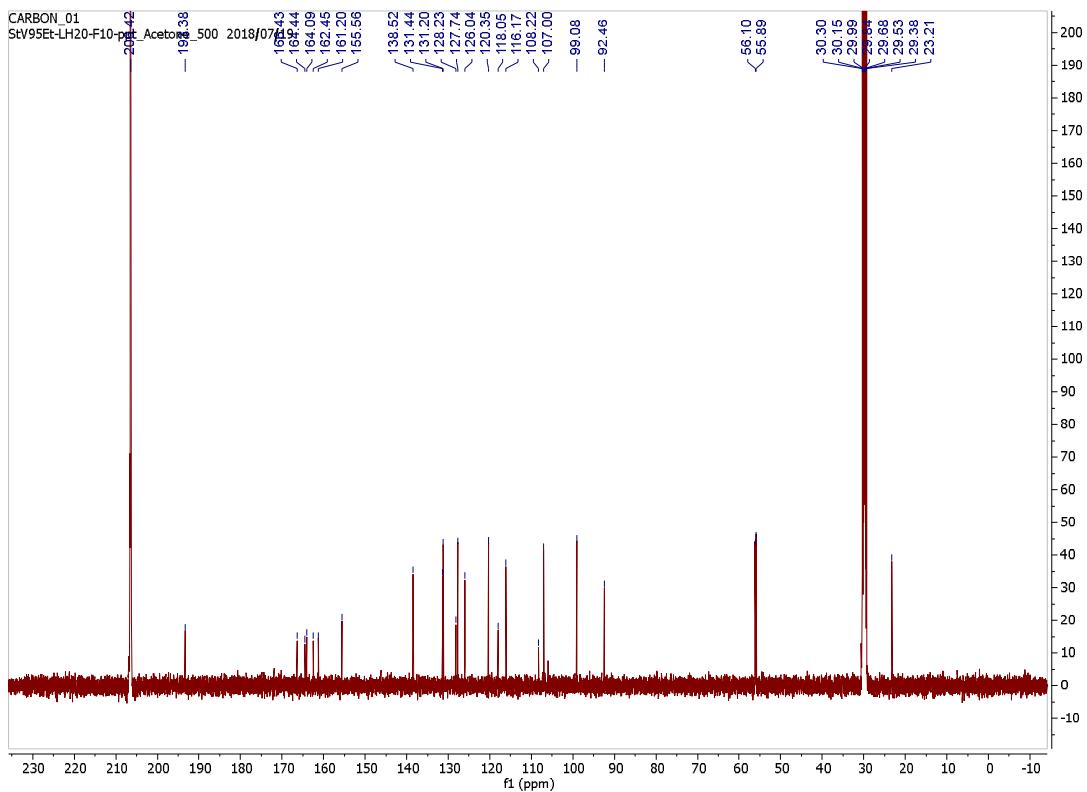


Figure S2. ¹³C NMR spectrum of ((E)-1-(3',5'-dihydroxy-4'-(2''-hydroxybenzyl)-2'-methoxyphenyl)-3-(2,4-dimethoxyphenyl) prop-2-en-1-one) (**1**) measured in 500 MHz in acetone

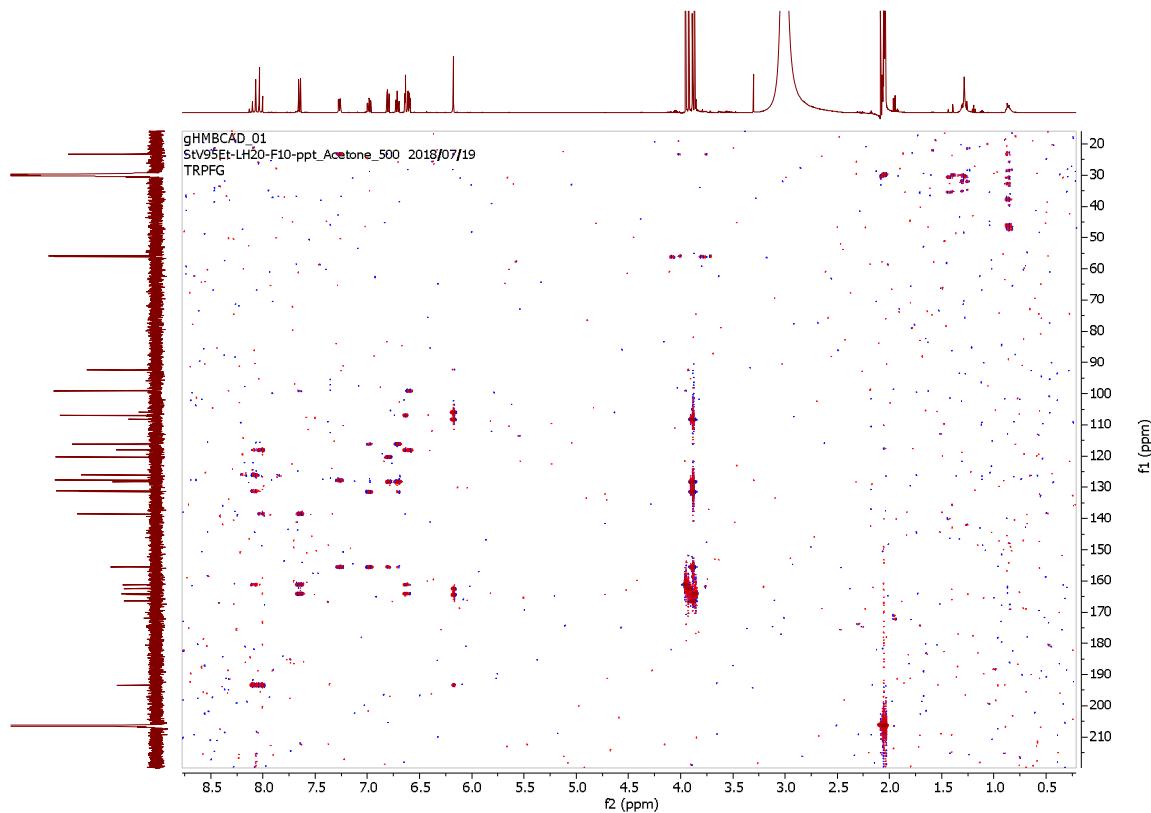


Figure S3. HMBC spectrum of ((E)-1-(3',5'-dihydroxy-4'-(2''-hydroxybenzyl)-2'-methoxyphenyl)-3-(2,4-dimethoxyphenyl) prop-2-en-1-one) (**1**) measured in 500 MHz in acetone

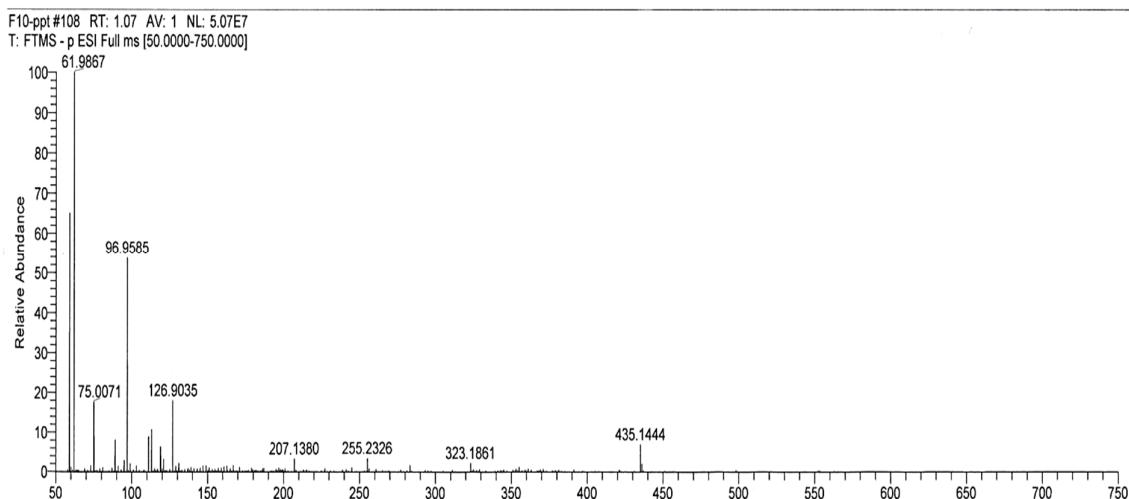


Figure S4. The HREIMS of (E)-1-(3',5'-dihydroxy-4'-(2''- hydroxybenzyl) -2'- methoxyphenyl) -3- (2,4-dimethoxyphenyl) prop-2-en-1-one (**1**), M= 436, [M - H]⁻

2. ¹H NMR and ¹³C-NMR spectrums, and HREIMS of (2E)-1-(2,4-Dihydroxy-6-methoxyphenyl)-3-(2,4-dimethoxyphenyl)-2-propen-1-one (**2**)

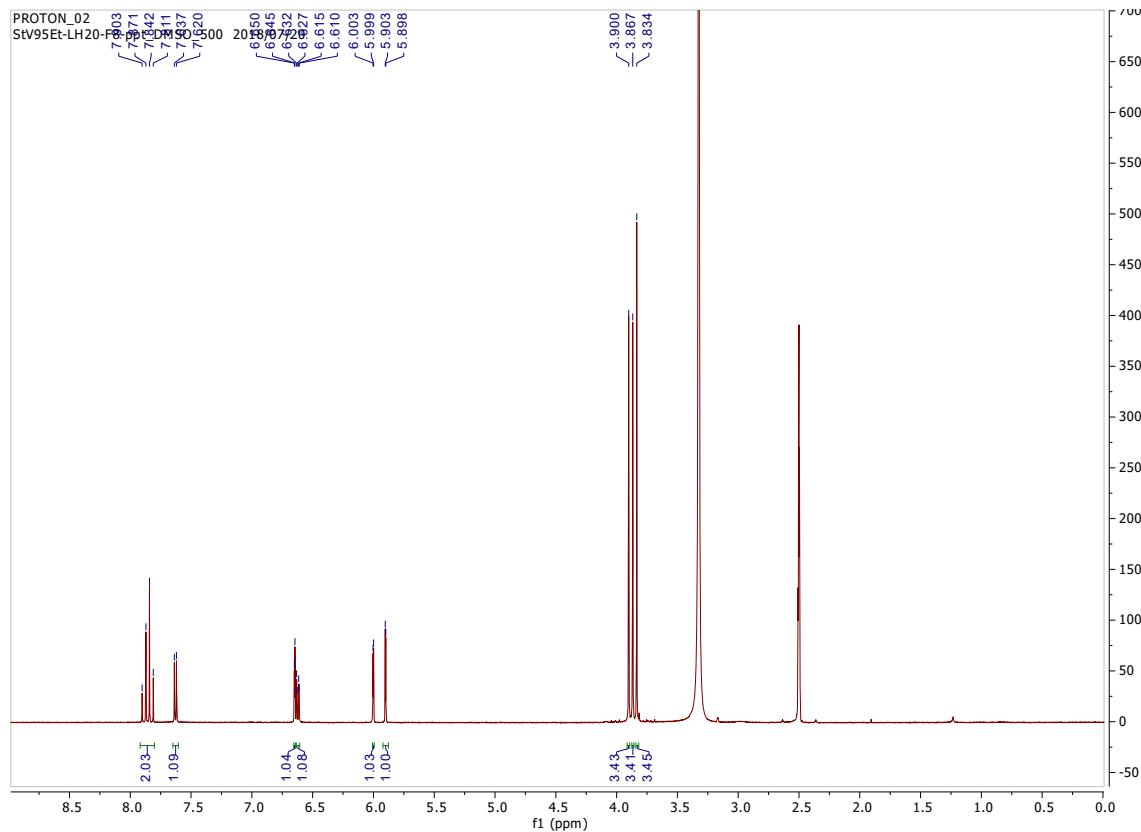


Figure S5. ¹H NMR spectrum of (2E)-1-(2,4-Dihydroxy-6-methoxyphenyl)-3-(2,4-dimethoxyphenyl)-2-propen-1-one (**2**) measured in 500 MHz in DMSO

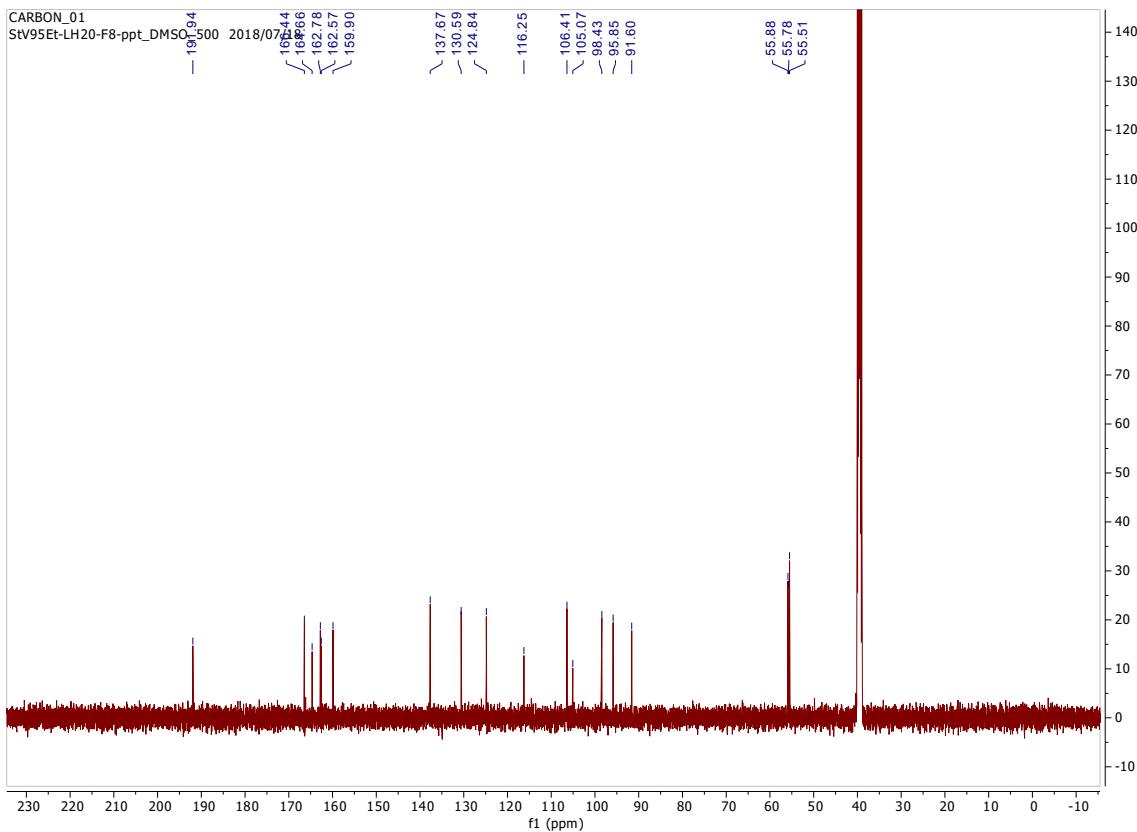


Figure S6. ^{13}C NMR spectrum of (2E)-1-(2,4-Dihydroxy-6-methoxyphenyl)-3-(2,4-dimethoxyphenyl)-2-propen-1-one (**2**) measured in 500 MHz in DMSO

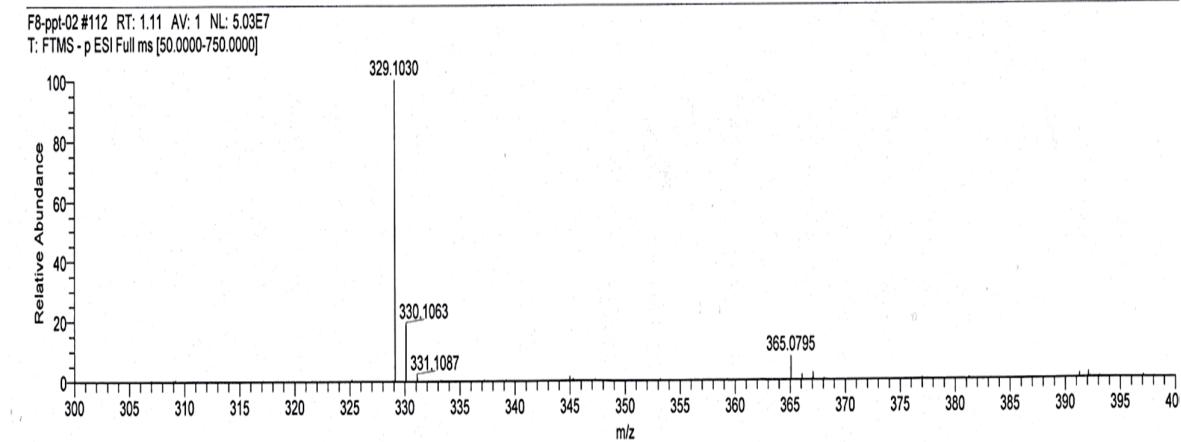


Figure S7. The HREIMS of (2E)-1-(2,4-Dihydroxy-6-methoxyphenyl)-3-(2,4-dimethoxyphenyl)-2-propen-1-one (**2**), M= 330, $[\text{M} - \text{H}]^-$

3. ^1H NMR and ^{13}C -NMR spectrums, and HREIMS of 7-hydroxyethyloxy-3',4',5-trimethoxy flavone (**3**)

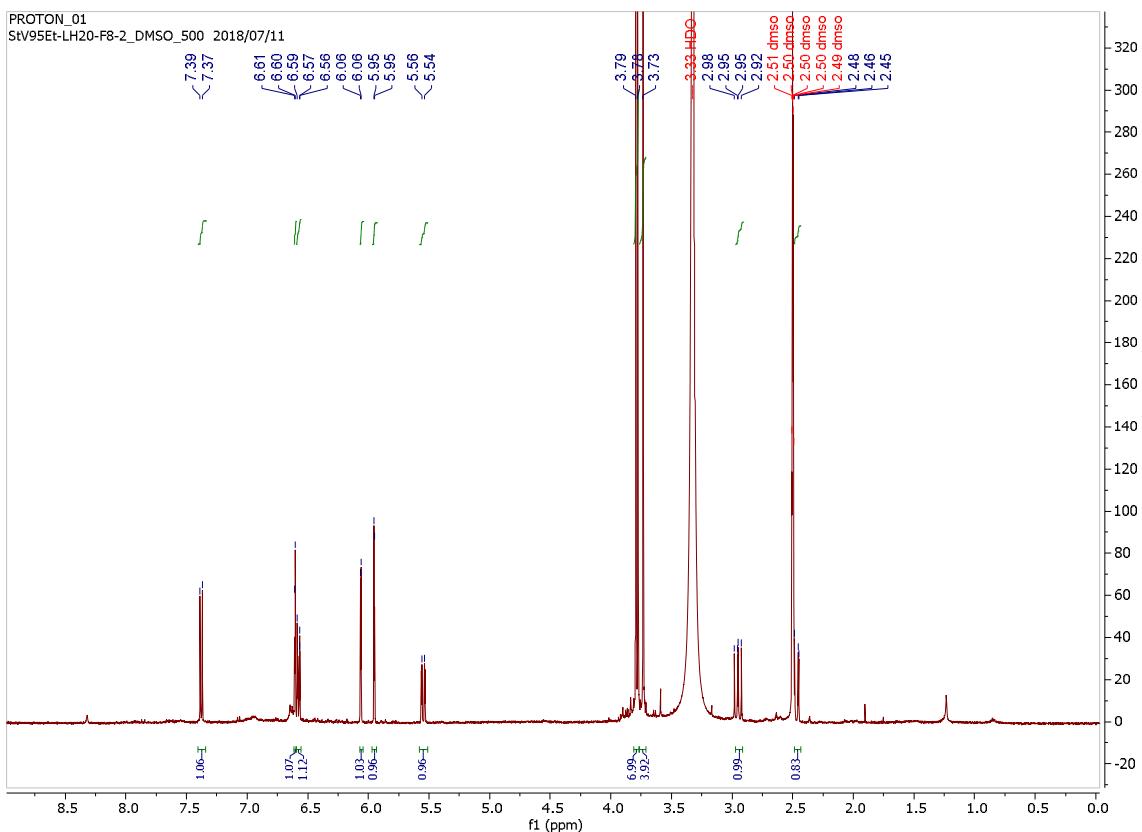


Figure S8. ^1H NMR spectrum of 7-hydroxyethoxy-3',4',5-trimethoxy flavone (**3**) measured in 500 MHz in DMSO

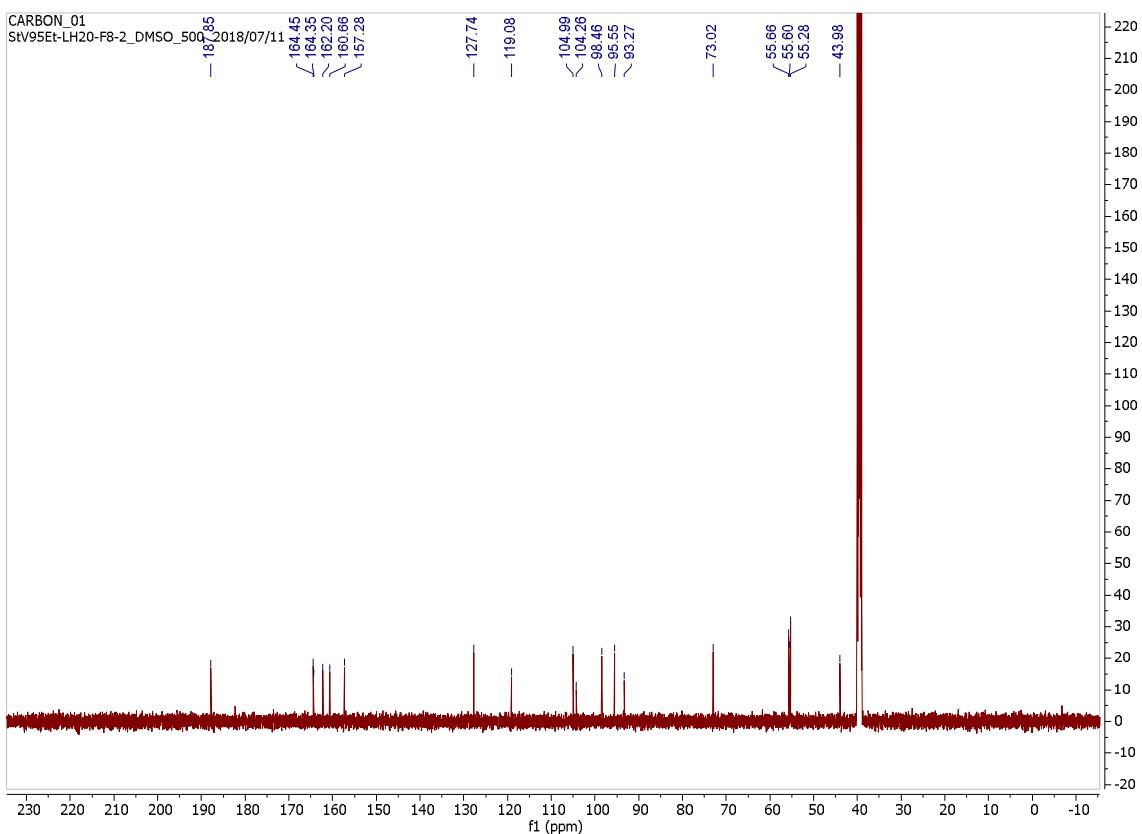


Figure S9. ^{13}C NMR spectrum of 7-hydroxyethoxy-3',4',5-trimethoxy flavone (**3**) measured in 500 MHz in DMSO

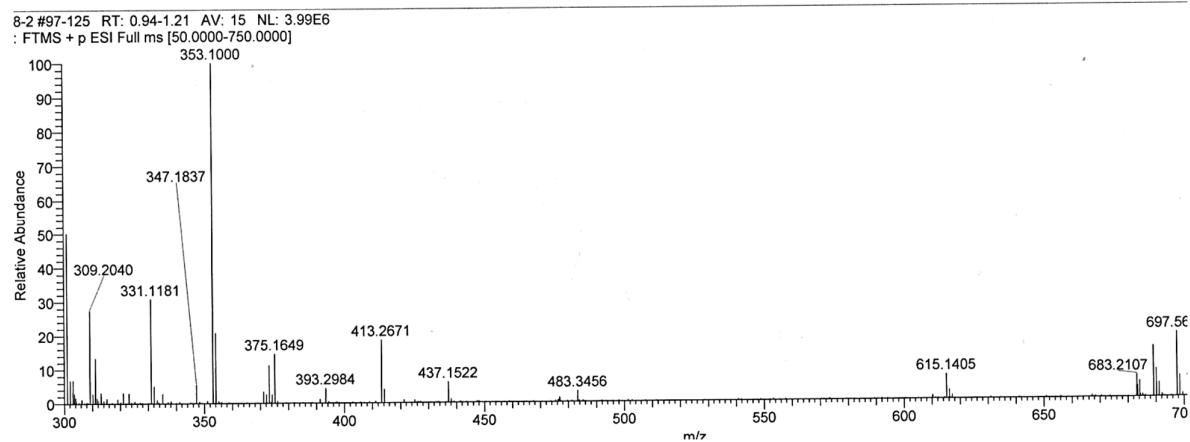


Figure S10. The HREIMS of 7-hydroxyethoxy-3',4',5-trimethoxy flavone (**3**), M= 330, [M + Na]⁺

4. ¹H NMR and ¹³C-NMR spectrums, and HREIMS of (+)-crotepoxide (**4**)

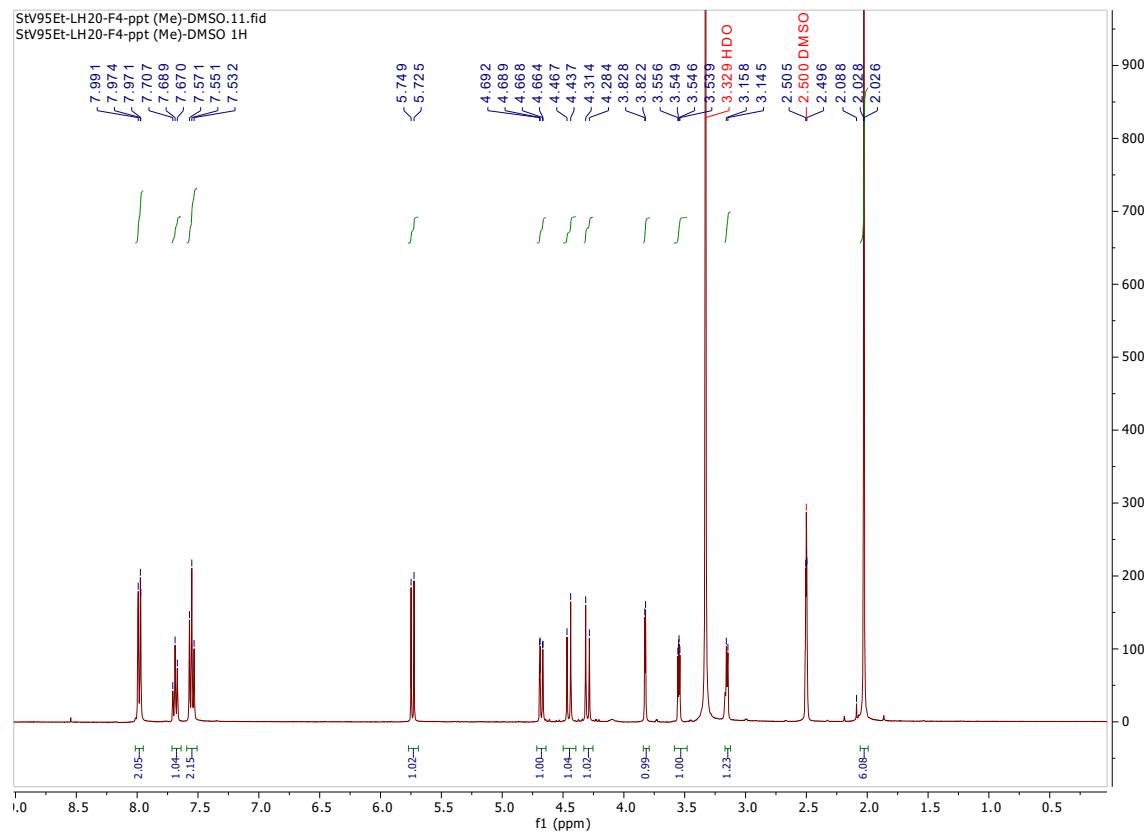


Figure S11. ¹H NMR spectrum of (+)-crotepoxide (**4**) measured in 500 MHz in DMSO

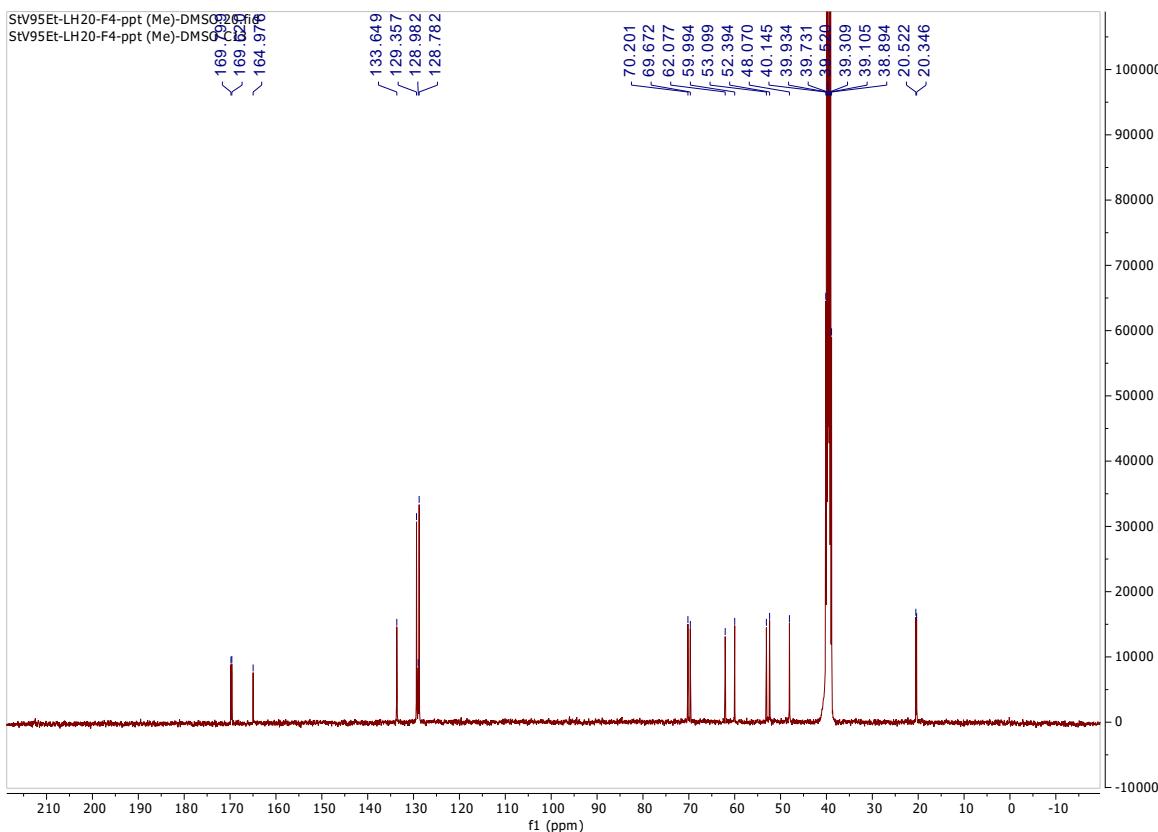


Figure S12. ^{13}C NMR spectrum of (+)-crotepoxide (**4**) measured in 500 MHz in DMSO

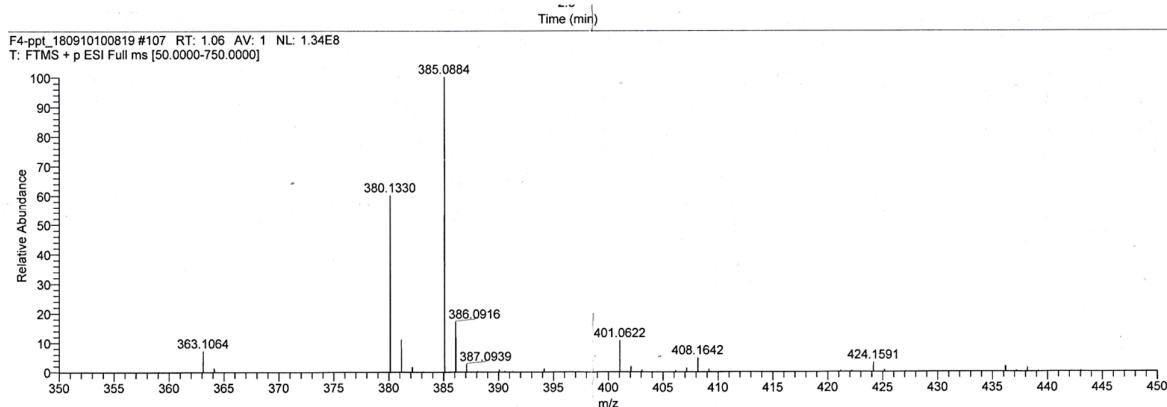


Figure S13. The HREIMS of (+)-crotepoxide (**4**), M= 362, $[\text{M} + \text{Na}]^+$

5. ^1H NMR and ^{13}C -NMR spectra, and HREIMS of sandaracopimamic acid (**5**)

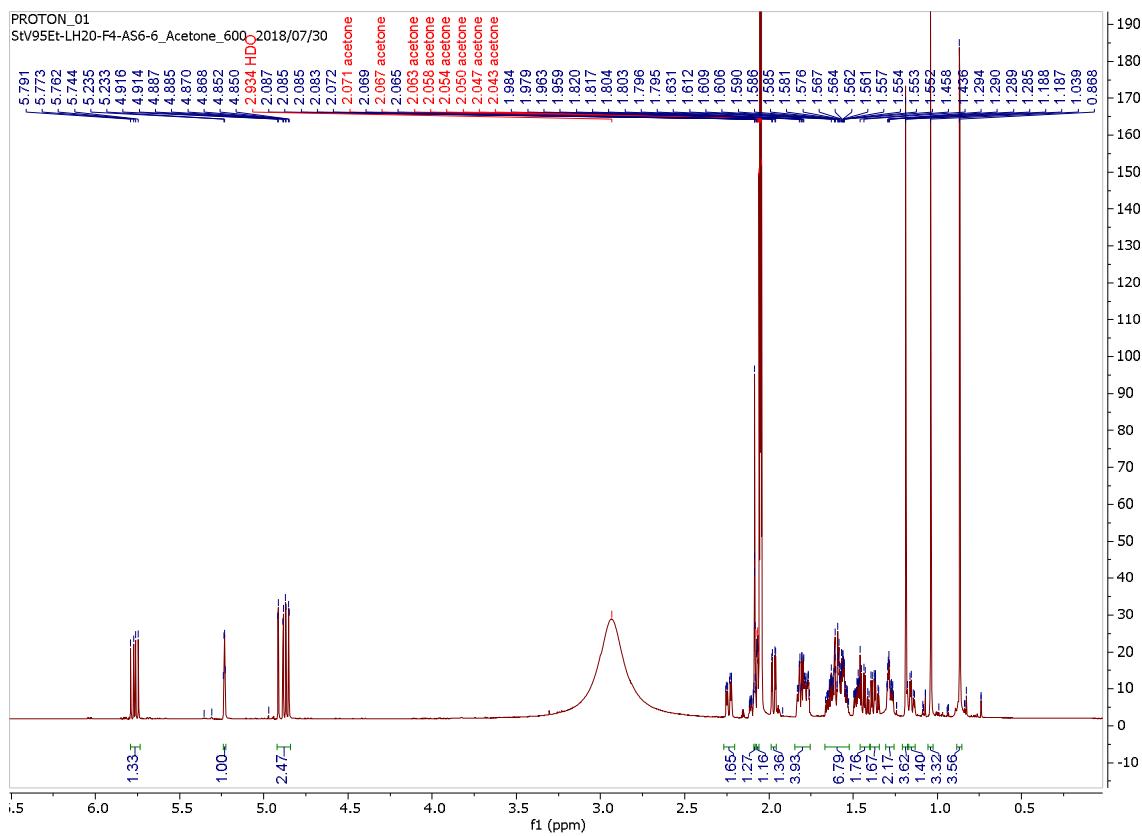


Figure S14. ^1H NMR spectrum of sandaracopimamic acid (**5**) measured in 500 MHz in acetone

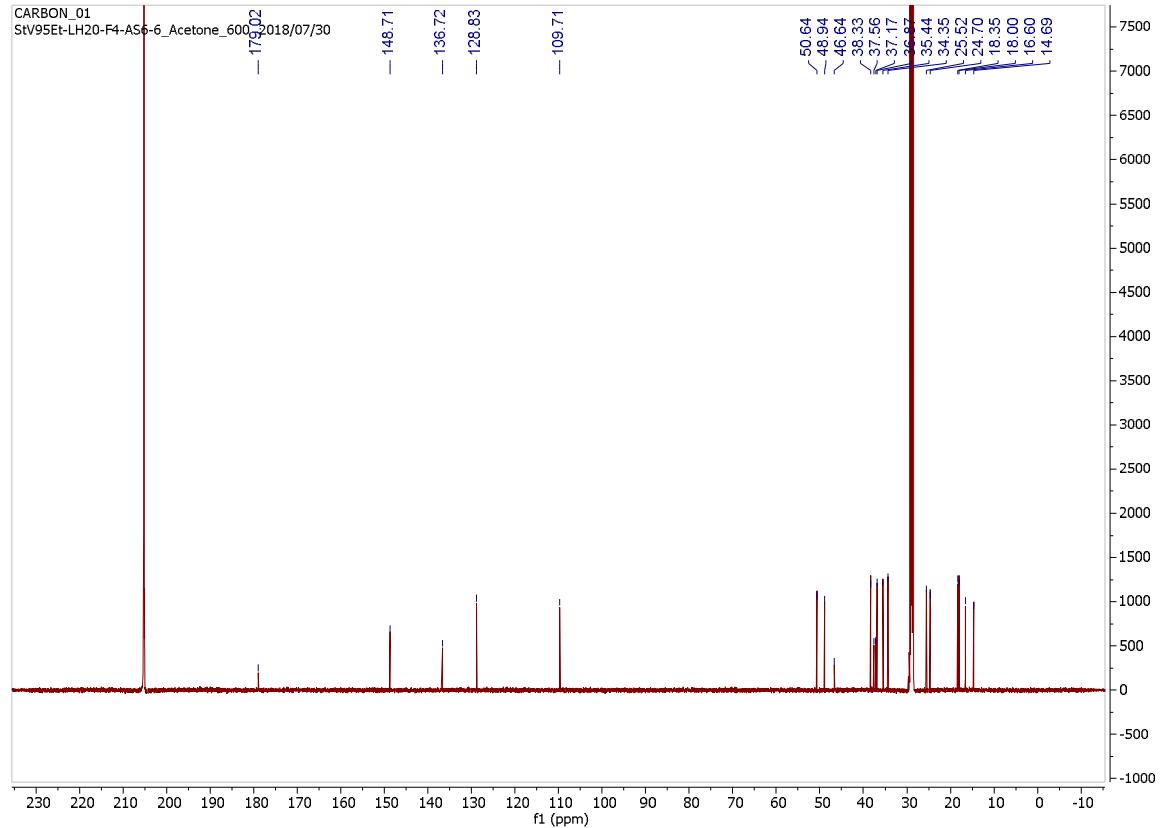


Figure S15. ^{13}C NMR spectrum of sandaracopimamic acid (**5**) measured in 500 MHz in acetone

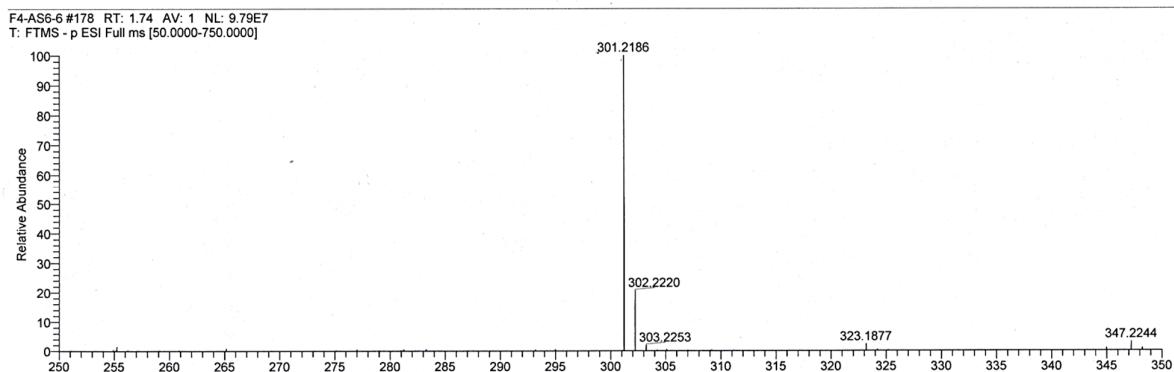


Figure S16. The HREIMS of sandaracopimamic acid (**5**), M= 302, [M - H]⁻

6. ^1H NMR and ^{13}C -NMR spectrums, and HREIMS of o-methoxybenzoyl benzoate (**6**)

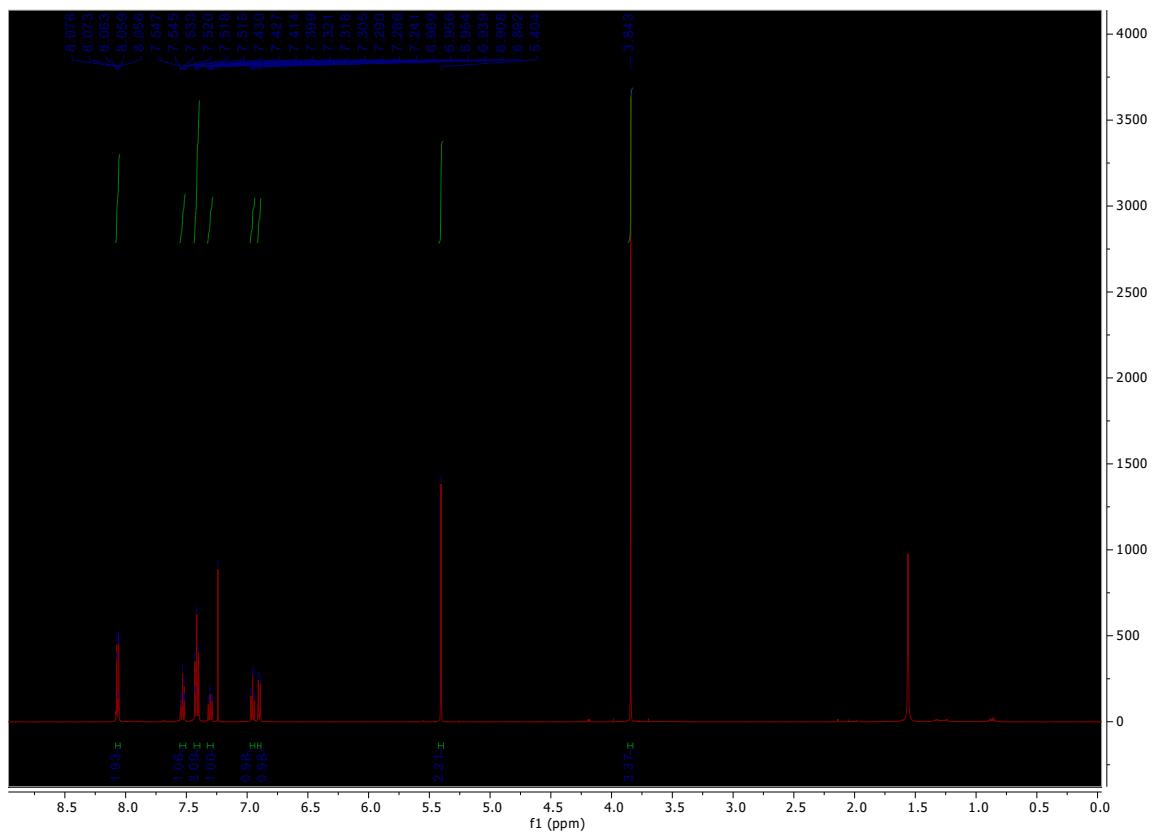


Figure S17. ^1H NMR spectrum of o-methoxybenzoyl benzoate (**6**) measured in 500 MHz in CDCl_3

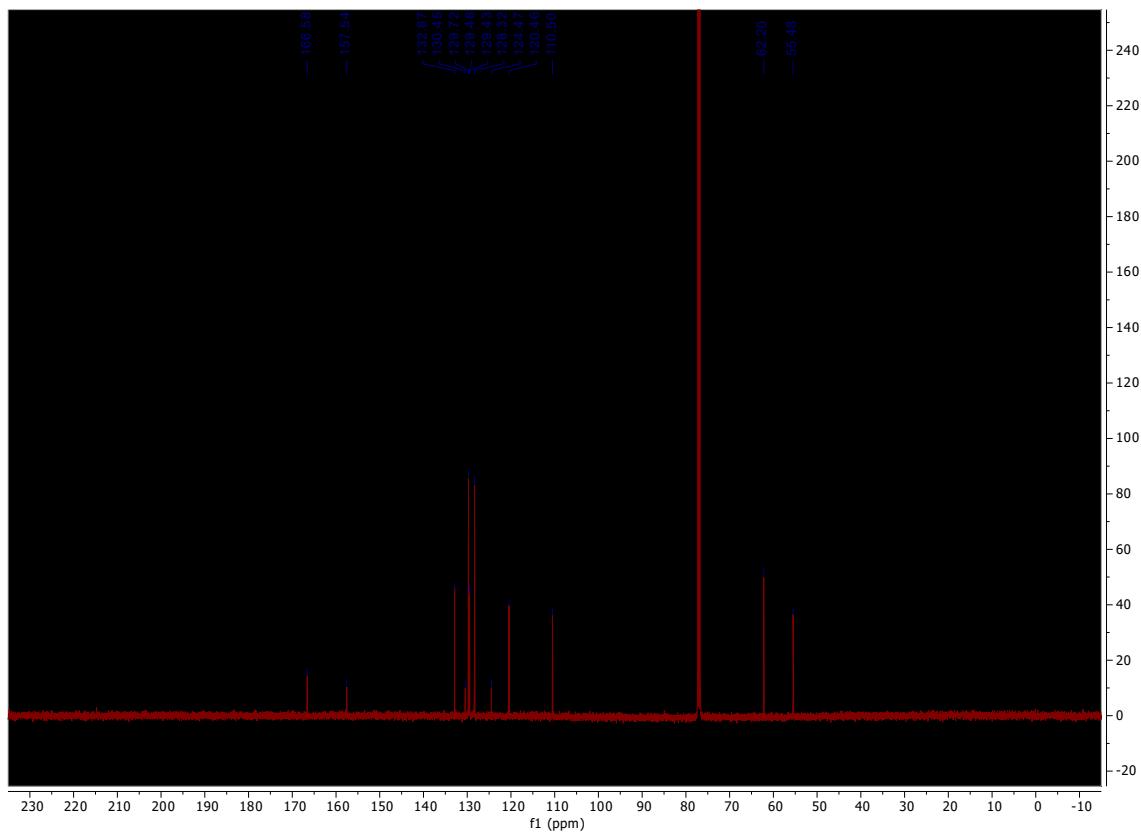


Figure S18. ^{13}C NMR spectrum of o-methoxybenzoyl benzoate (**6**) measured in 500 MHz in CDCl_3

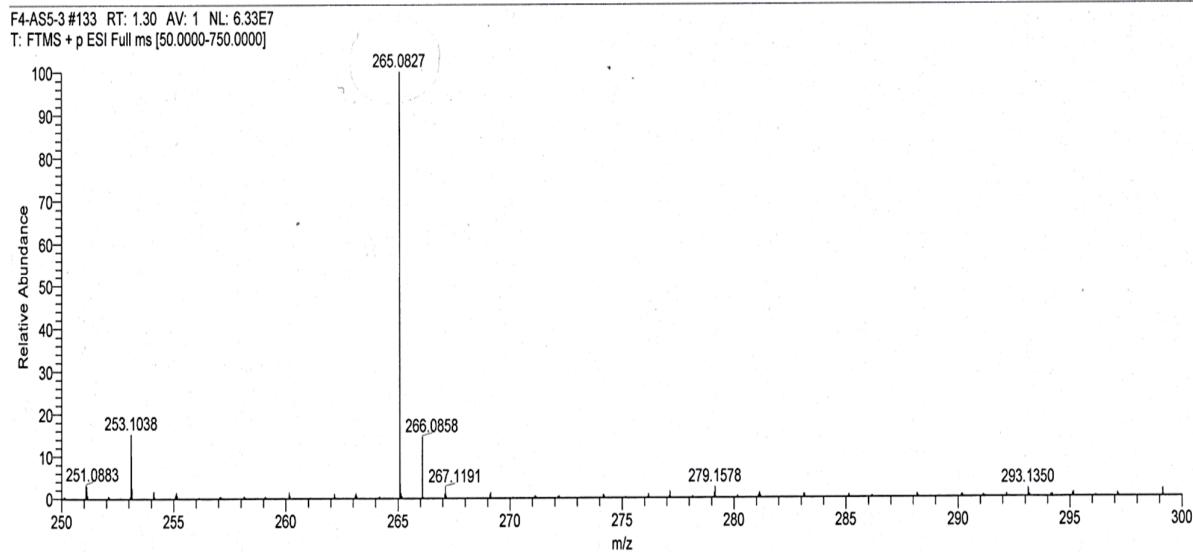


Figure S19. The HREIMS of o-methoxybenzoyl benzoate (**6**), $M = 242 [\text{M} + \text{Na}]^+$

7. ^1H NMR and ^{13}C -NMR spectrums, and HREIMS of (-)-1,6-desoxytingtanoxide (**7**)

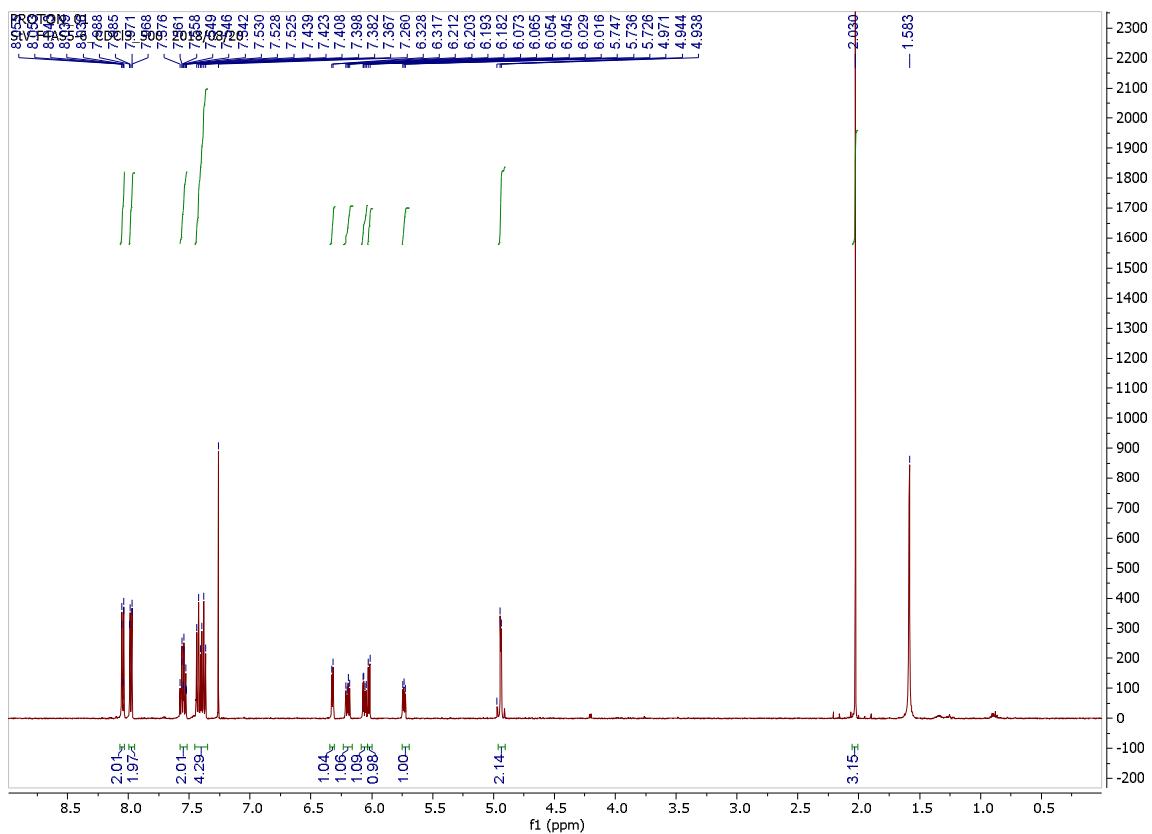


Figure S20. ^1H NMR spectrum of (-)-1,6-desoxytingtanoxide (7) measured in 500 MHz in CDCl_3

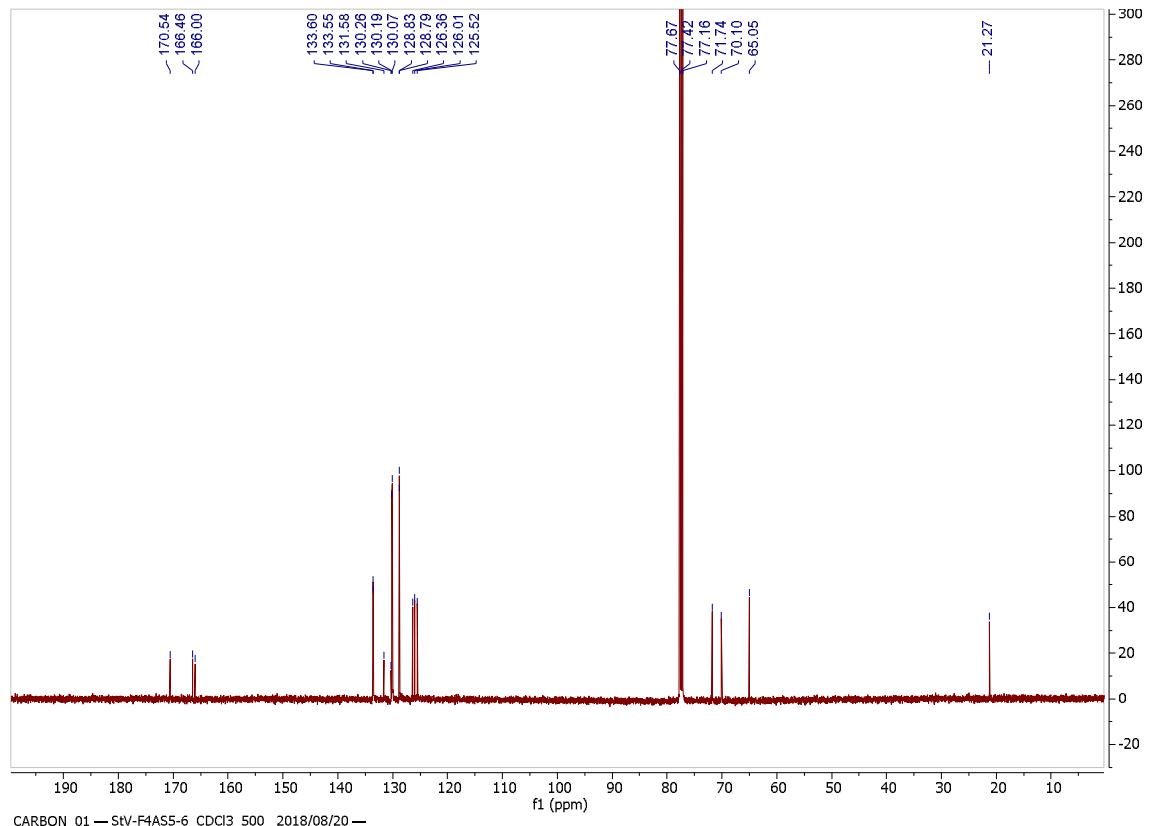


Figure S21. ^{13}C NMR spectrum of ($-$)1,6-desoxytingtanoxide (**7**) measured in 500 MHz in CDCl_3

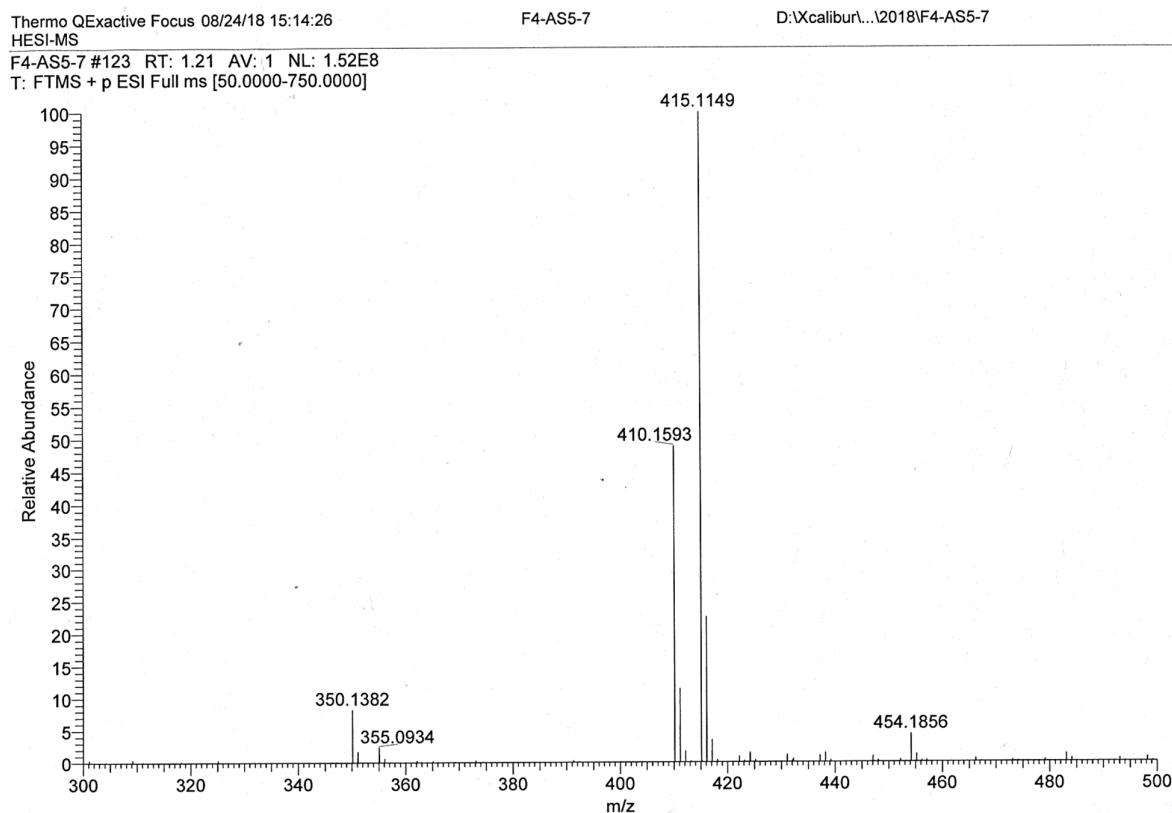


Figure S22. The HREIMS of ($-$)1,6-desoxytingtanoxide (**7**), $M= 392$, $[\text{M} + \text{Na}]^+$