

SUPPLEMENTARY INFORMATION

Supplementary Table 1

Calibration of radiocarbon data on ancient *Commiphora* seed using OxCal 4.4 (1-3)

ETH ^a no.	UZ ^b no	C-14 age ^c	$\delta^{13}\text{C}$ (‰)	Calibration	
				1 σ (68.3%)	2 σ (95.4%)
40349	5834	940 ± 35	-18.1 ± 1.0	1040 – 1156 cal CE	1026 – 1202 cal CE

^a Swiss Federal Institute of Technology, ^b UZ (University of Zurich)

^c Sample (fraction of the operculum) weight was 8.8 mg giving a CO₂ pressure in the reactor of 503 mbar. Organic C in the cracked sample was 41.0%. One sample was analysed (no repeat available due to small sample size). All data produced is shown here.

Supplementary Table 2

Measured and modelled (using different contamination scenarios) C-14 ages of the seed shell.

HU LMK G1.SF (sample name)				Calibrated ages (2 s)	Median age	
ETH no. 40349	UZ no 5834	C-14age	SD	pMC (%)	calCE	(calCE)
measured	940	35	88.3	1026 -	1202	1103
1% addition of fresh C*	957	35	88.1	1021 -	1169	1099
2% addition of fresh C*	972	35	87.9	996 - 1161	1095	
3% addition of fresh C*	988	35	87.8	993 - 1157	1085	

* Corrected radiocarbon age assuming that 1, 2 or 3% of modern carbon was absorbed in the sample.

Supplementary Table 3

List of major compounds detected in “Sheba” resin by GC-MS (indicated in Supplementary Fig 2)

Peak	RT	Compound
1	8.47	α -pinene*
2	9.52	β -pinene*
3	10.69	limonene*
4	42.89	unknown triterpene C ₃₀ H ₅₀ O
5	42.98	β -amyrin*
6	43.34	unknown triterpene C ₃₀ H ₅₀ O
7	43.48	lupeol*
8	43.52	α -amyrin*

*confirmed with commercial standard

Supplementary Table 4

¹H (500 MHz, in CDCl₃) and ¹³C (125 MHz) chemical shifts for compound II (cf Fig 4 for structure and Supplementary Fig 5 for structure and carbon numbering)

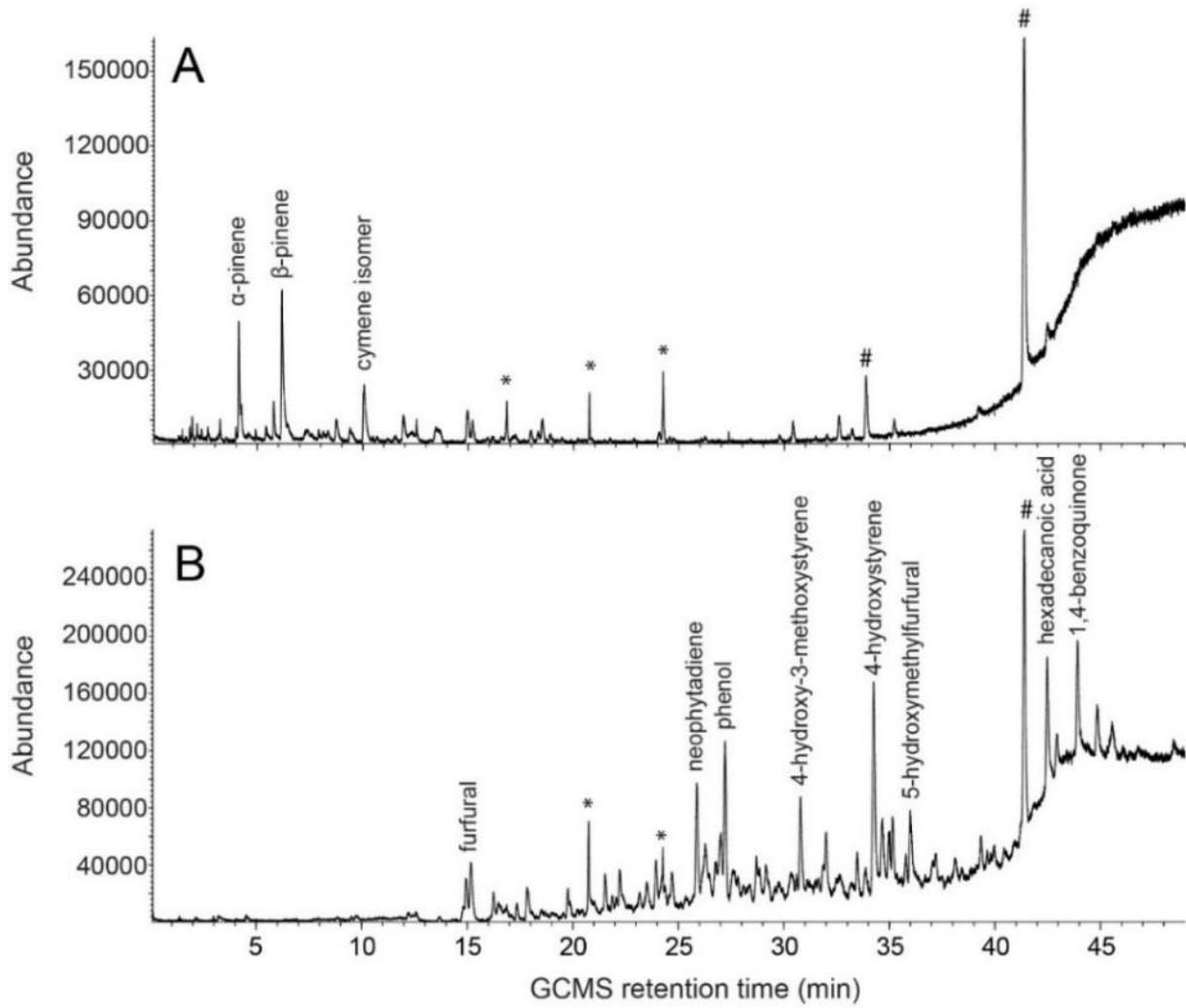
Position	δ_c	δ_h	δ_h	Position	δ_c	δ_h	δ_h
I ₁	62.1	4.30	4.16	X ₁	101.2	4.63	
I ₂	71.8	5.14		X ₂	70.6	4.84	
I ₃	71.2	5.09		X ₃	71.8	5.09	
I ₄	25.6	1.62	1.28	X ₄	69.0	4.90	
I ₅	28.9	1.50	1.50	X ₅	61.9	4.10	3.38
I ₆	78.5	3.58		r ₁	96.8	4.76	
I ₇	34.5	1.54	1.44	r ₂	72.3	5.05	
I ₈	29.5	1.26	1.26	r ₃	74.8	4.00	
I ₉₋₁₇	29.8	1.26	1.26	r ₄	72.8	5.03	
I ₁₈	32.1	1.25		r ₅	66.8	3.81	
I ₁₉	22.8	1.29		r ₆	17.4	1.150	
I ₂₀	14.3	0.877		Acetates (CO) 169.4-170.8			
				Acetates (Me) 20.7-21.1		1.98-2.14	

Supplementary Table 5

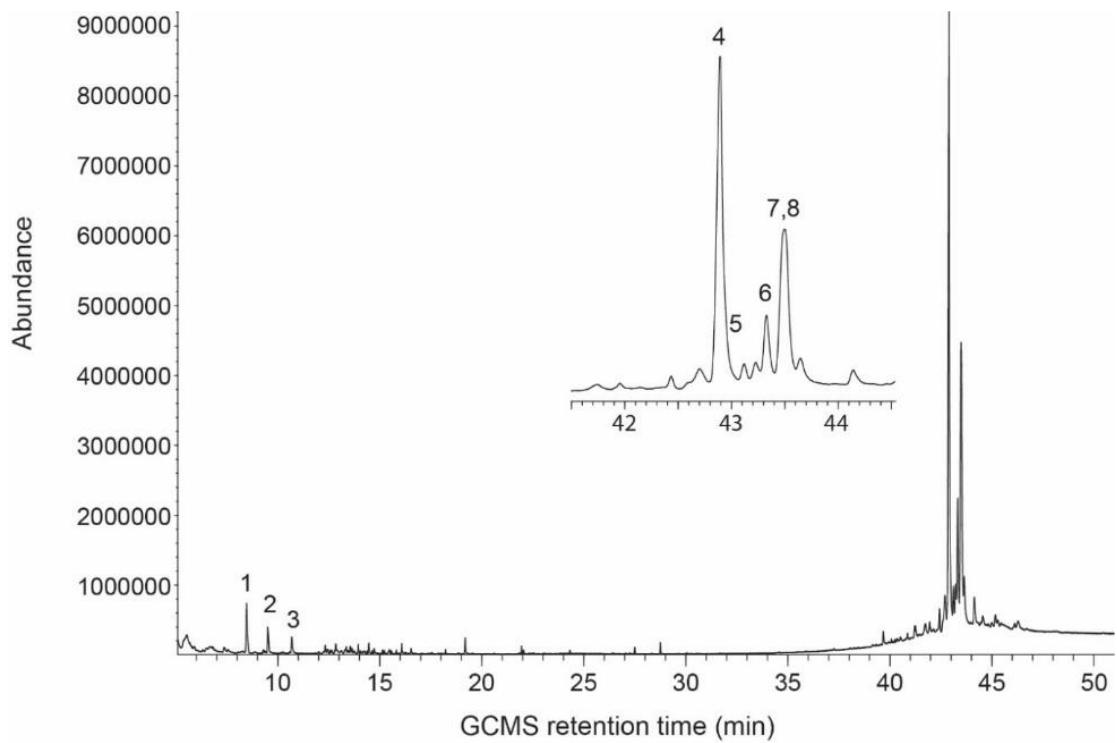
Authors and sources in antiquity discussing Judean Balsam

Author	Period	Source
Theophrastus	4 th century BCE	<i>Hist. Plant.</i> 9: 6:1–4; 9:7:3
Diodorus Siculus,	1 st century BCE	<i>Bibl. Hist.</i> 2 :48:9; 19:98
Strabo	1 st century BCE -1 st century CE	<i>Geogr.</i> 16:2:41; 17:1:15 and 16:4:19)
Gnaeus Pompeius Trogus (excerpted by Justin in the 2 nd century CE)	1 st century BCE	<i>Epitoma Historiarum Philippicarum</i> 36:3:1–5 https://www.attalus.org/translate/justin5.html
Dioscorides Pedanius	1 st century CE	<i>Mat. Med.</i> 1:19:1
Josephus	1 st century CE	<i>Flavius Josephus, War</i> 1:138, 361, <i>Ant.</i> 4:100; 14:54; 15:96, <i>Ant.</i> 8: 6: 6.
Pliny the Elder	1 st century CE	<i>Historia Naturalis</i> 12: 54 (25). Ed. Bostock J, (Taylor and Francis, London 1855)
Tacitus	1 st century CE	<i>Tacitus, Hist.</i> 5: 6: 1. (Loeb Classical Library Edition Tacitus https://penelope.uchicago.edu/Thayer/E/Roman/Texts/Tacitus/home.html
Galen	2 nd century CE	<i>De Antidotis</i> 1.4
Solanus	3 rd century CE	<i>Collectanea</i> 35.5.5-6
Eusebius of Caesarea (Eusebius Pamphili)	4 th century CE	<i>Onomasticon</i> 42:1-5:
Saint Jerome, Hieronymus	4-5 th century CE	<i>Letter 108 to Eustochium</i> , 11. 5
Bede (Beda Venerabilis)	7 th -8 th century CE	<i>Loc. Sanct.</i> 9.3/313
St Willibald	8 th century CE	<i>Hodeporicon</i> : Transl: Rev Canon Brownlow, (Palestine Pilgrims Text Society 1891)
Talmud *	3 rd -5 th century CE	Jerusalem Talmud: <i>Yoma</i> 4:5; Tractate Shabbat 25b, 26a
Talmud *	3 rd -6 th century CE	Babylonian Talmud <i>Kritut</i> 6a,

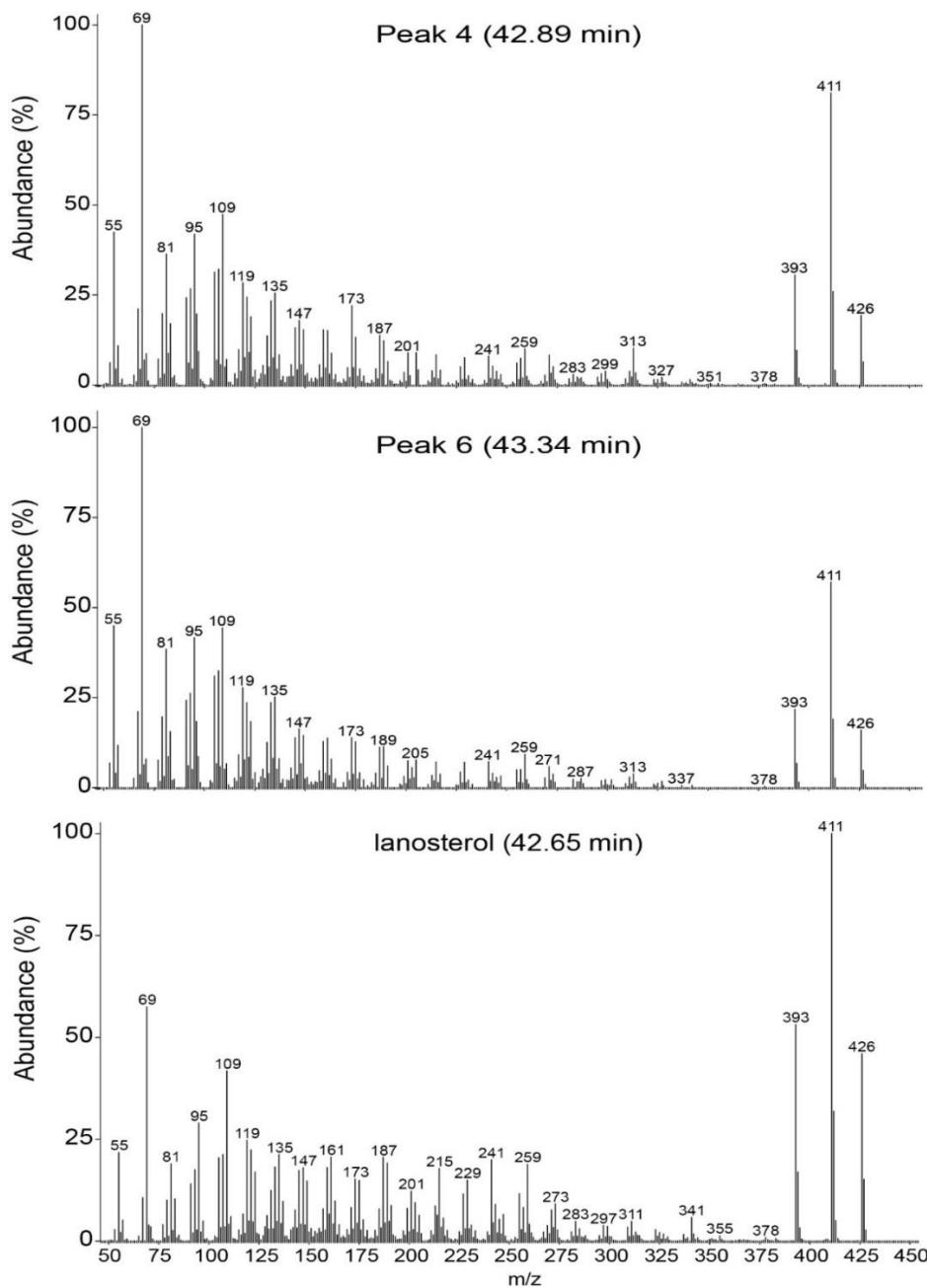
* Referring to Biblical “*tsori* ”



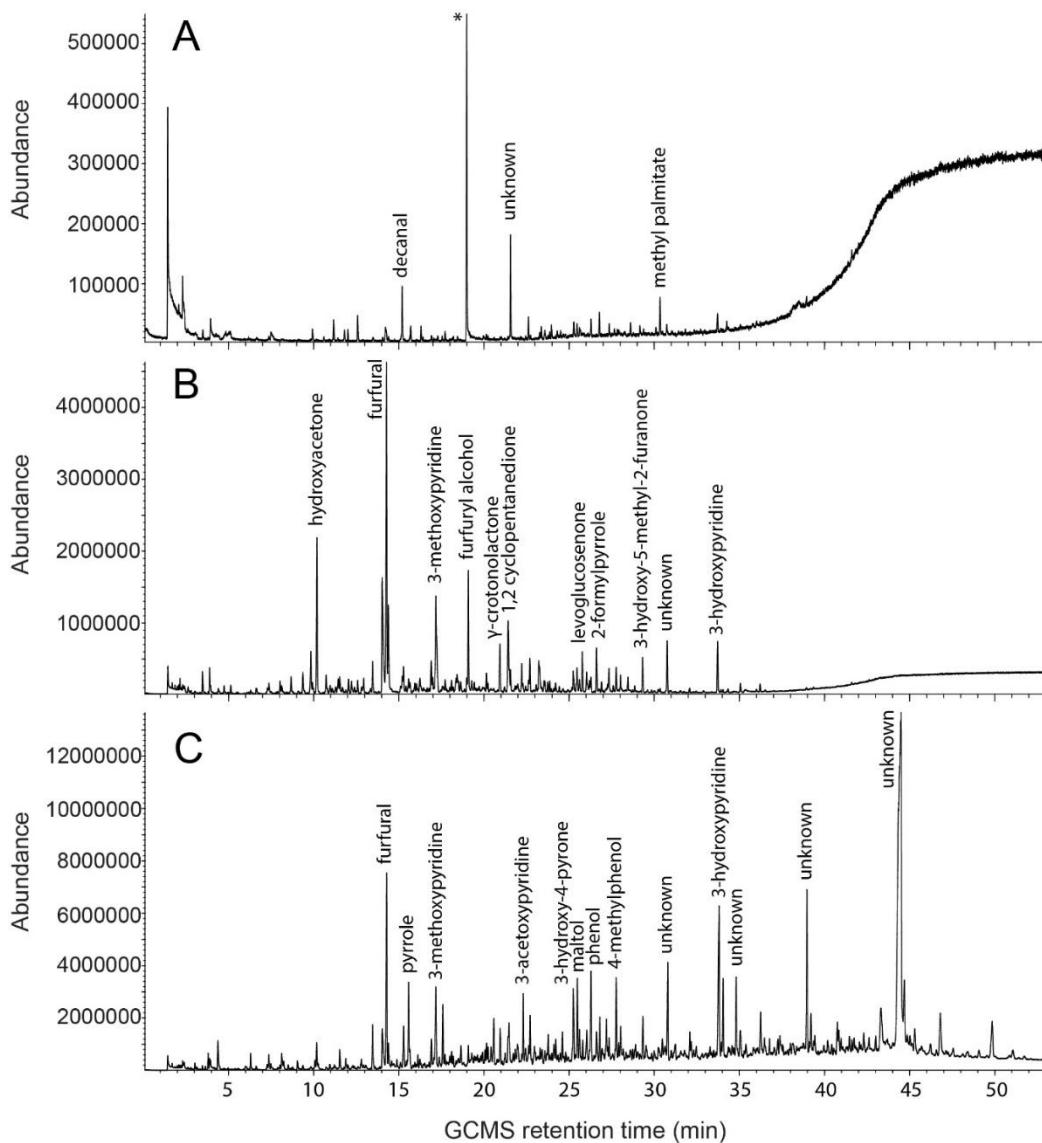
Supplementary Fig 1: Comparison of volatile compounds by SPME/GC-MS analysis from “Sheba” plant material (leaves and stems).



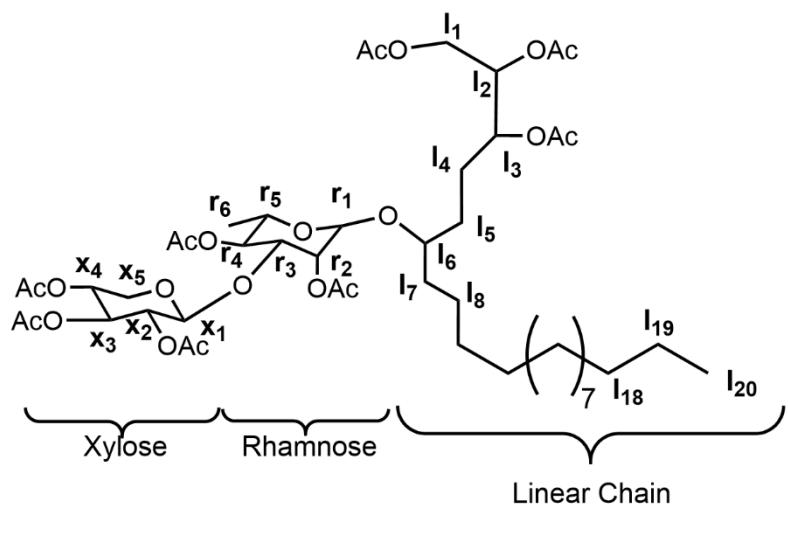
Supplementary Fig 2: Analysis of the hexanes extract of “Sheba” resin by GC-MS. Peaks marked with numbers correspond with identified compounds in **Supplementary Table 2**.



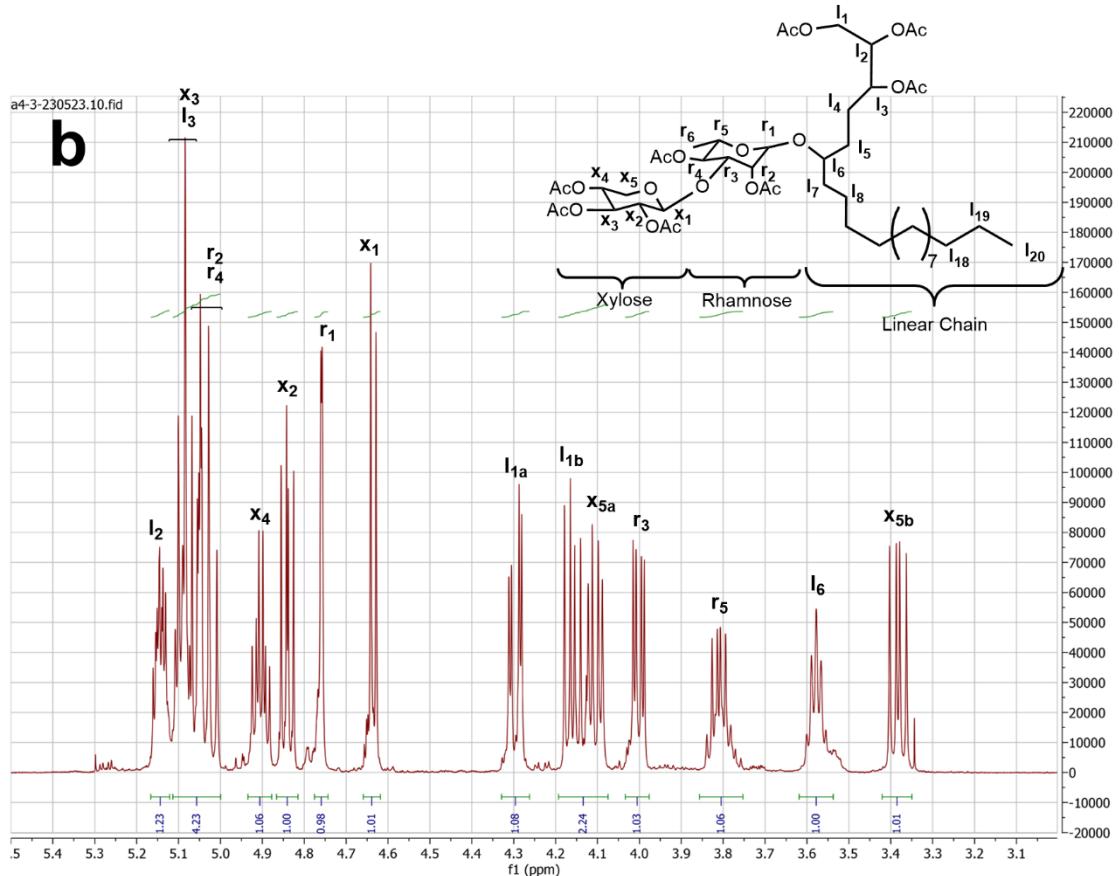
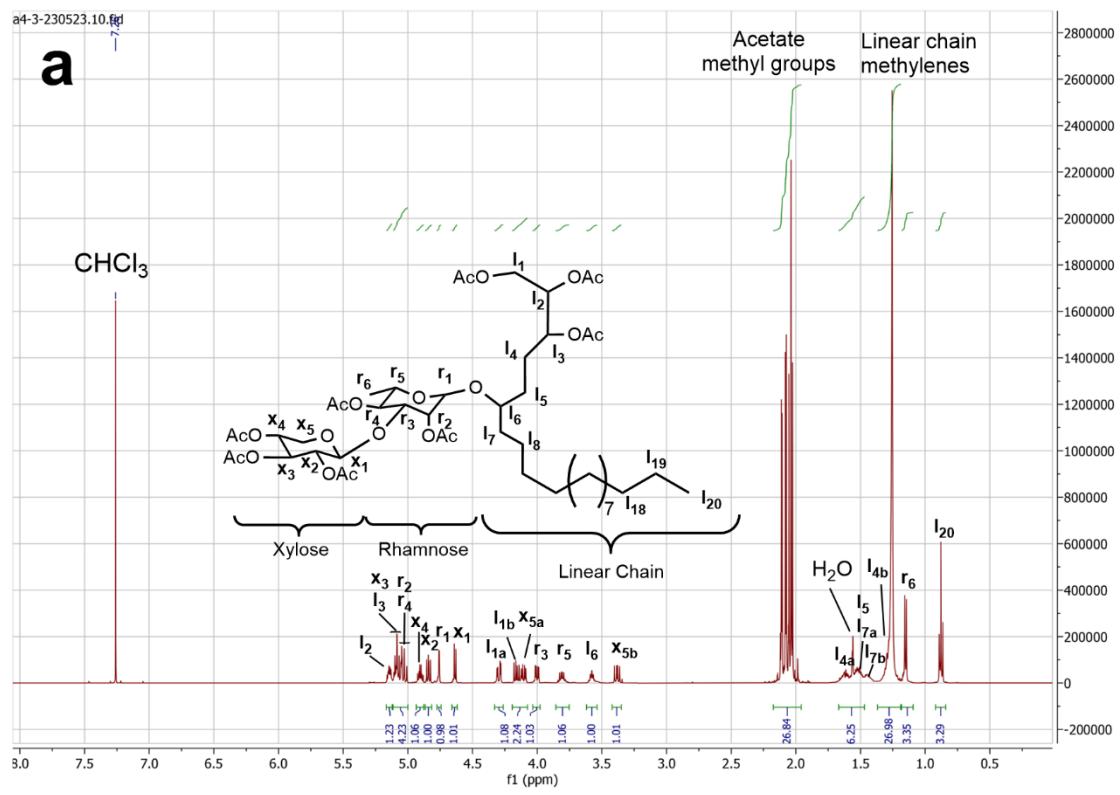
Supplementary Fig. 3: Mass spectra of the two unknown triterpenoids from “Sheba” resin (corresponding to **Supplementary Fig 2** and **Supplementary Table 2**) compared with lanosterol analysed under identical GC-MS conditions.



Supplementary Fig 4: Comparison of volatiles emitted from heating “Sheba” resin for increasing time and resulting volatile compounds trapped by SPME. **A.** 1 min, low heat. **B** 2 min, medium heat. **C** 5 min, high heat. *siloxane impurity.

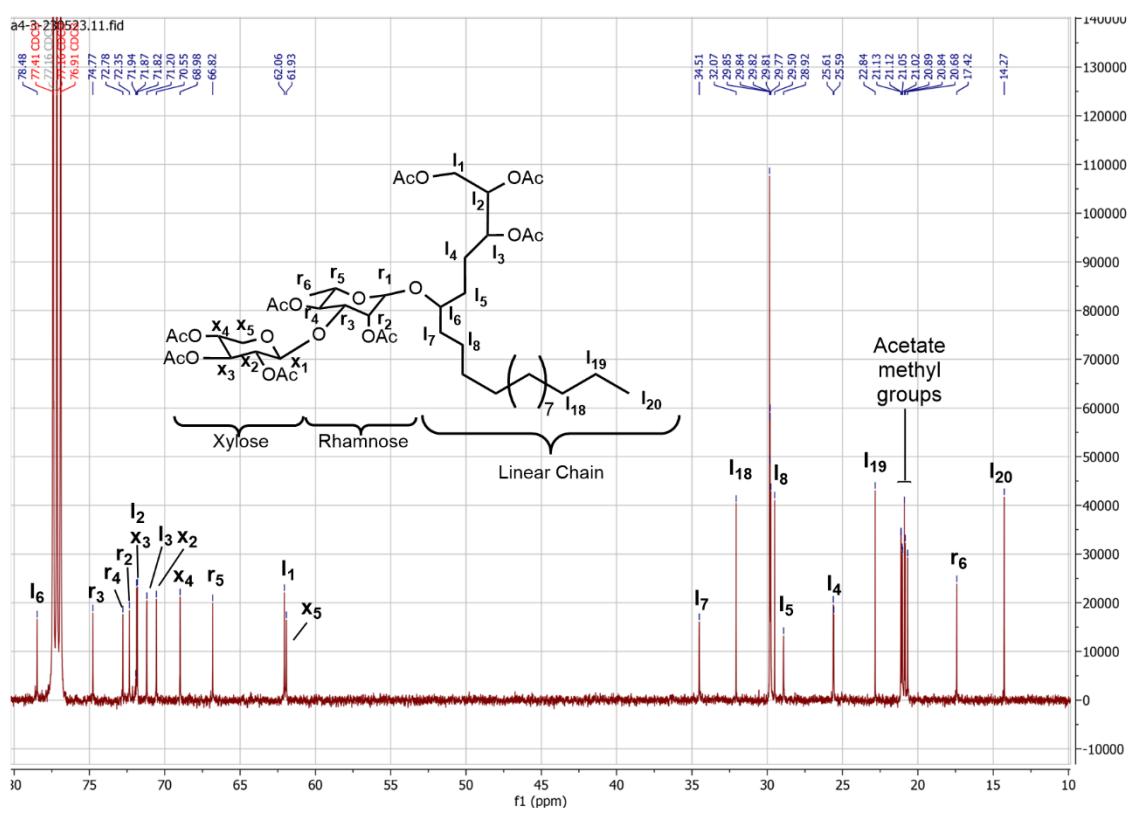
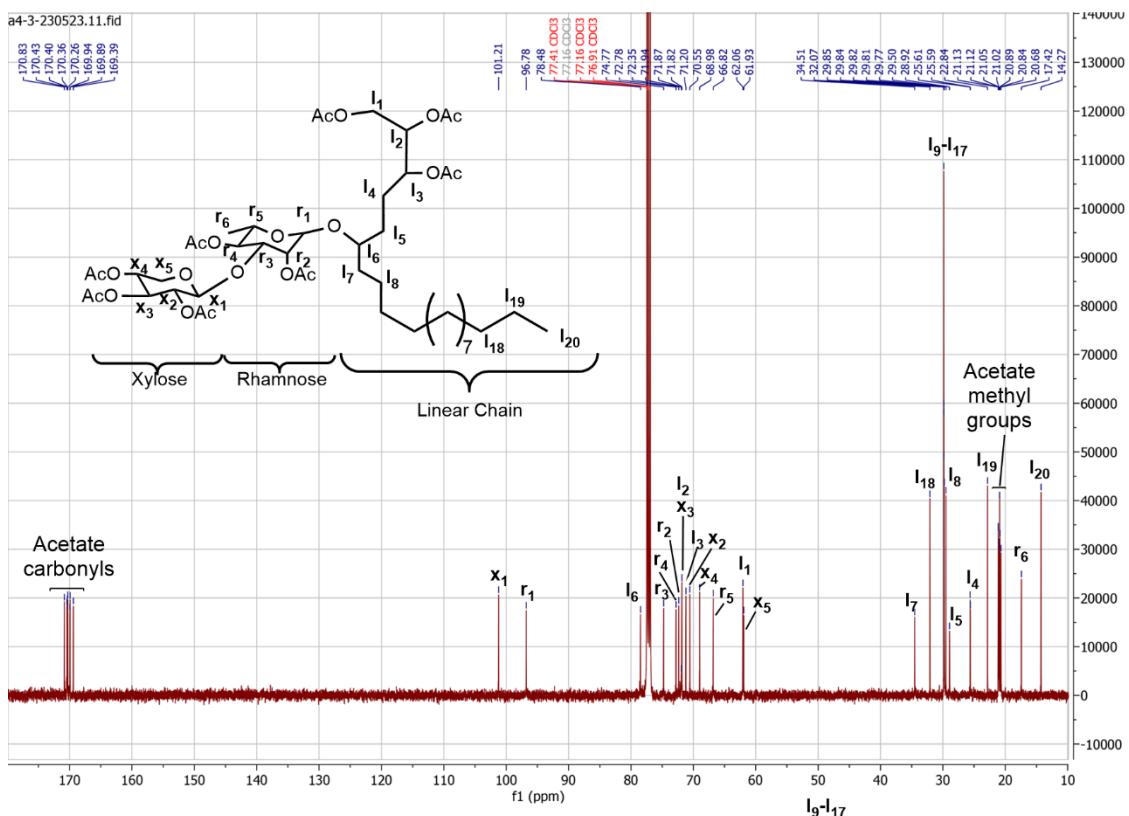


Supplementary Fig 5: Structure of compound **II** isolated from “Sheba” resin (cf **Fig. 4**) and identified using Nuclear Magnetic Resonance and High Resolution Mass Spectrometry.



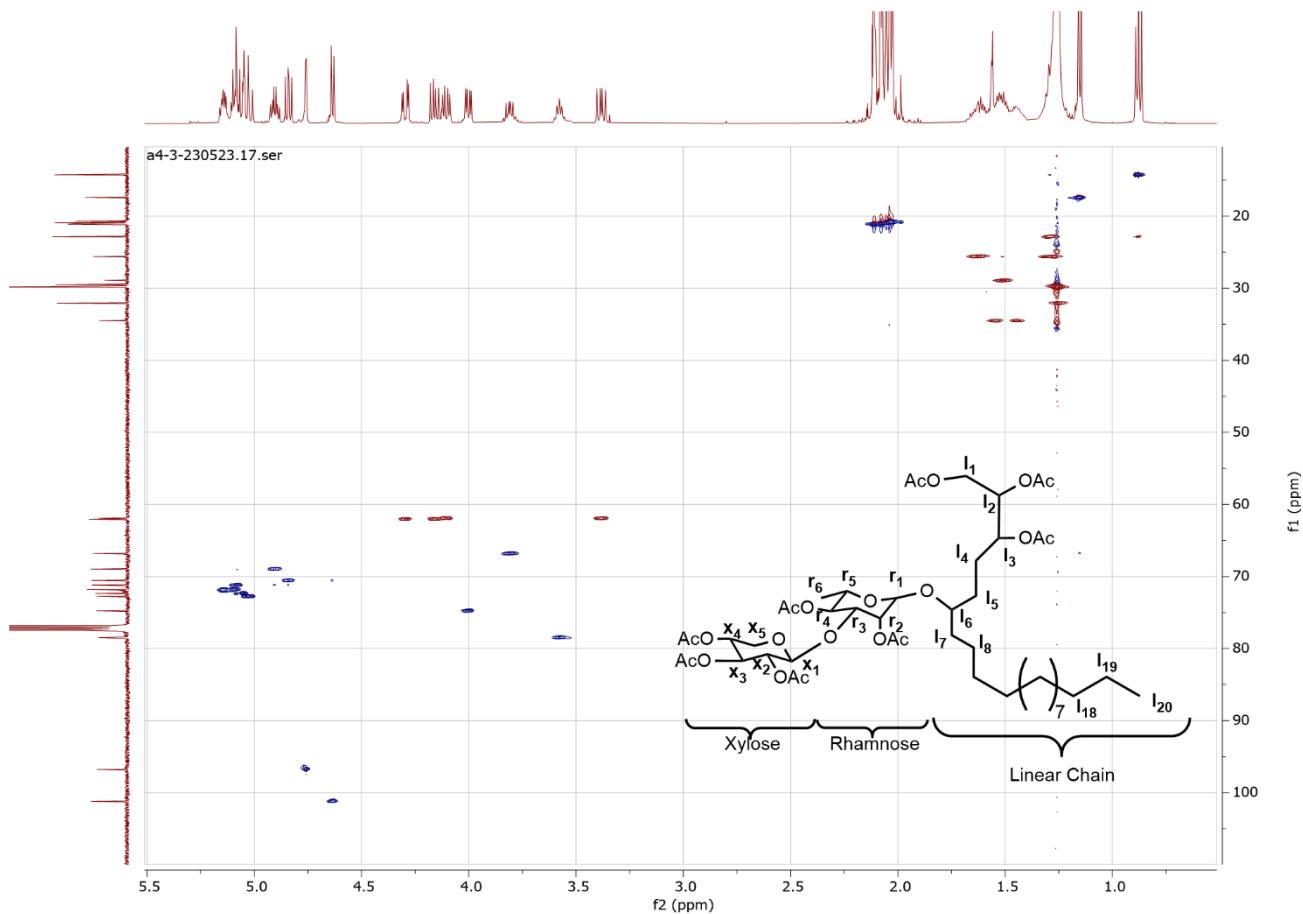
Supplementary Fig 6: ¹H-NMR spectrum (500 MHz, CDCl₃) of compound **II** (a) 0.0-8.0 ppm range;

(b) 3.0-5.5 ppm range.

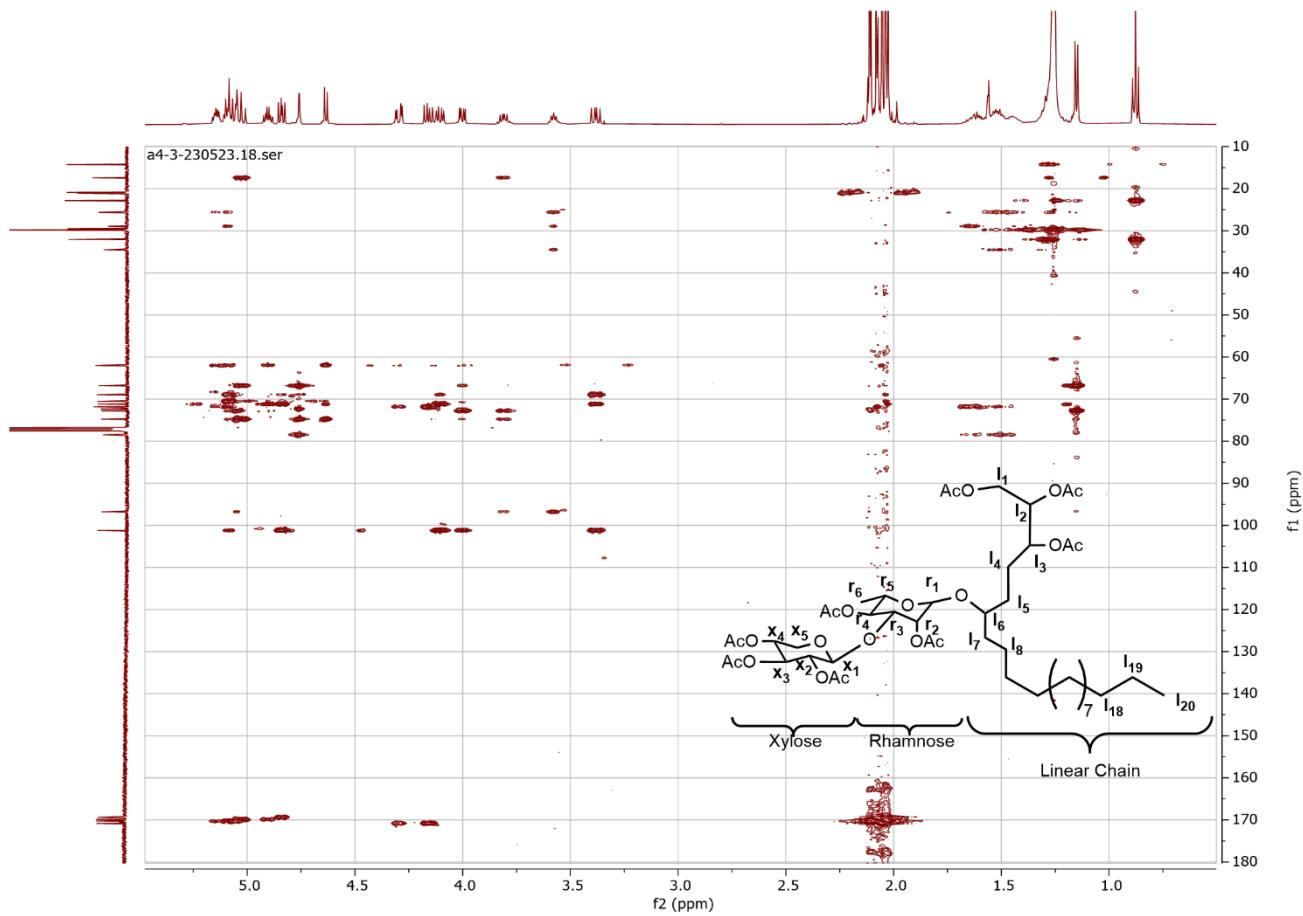


Supplementary Fig 7: ^{13}C -NMR spectrum (125 MHz, CDCl_3) of compound **II (a)** 10-190 ppm range;

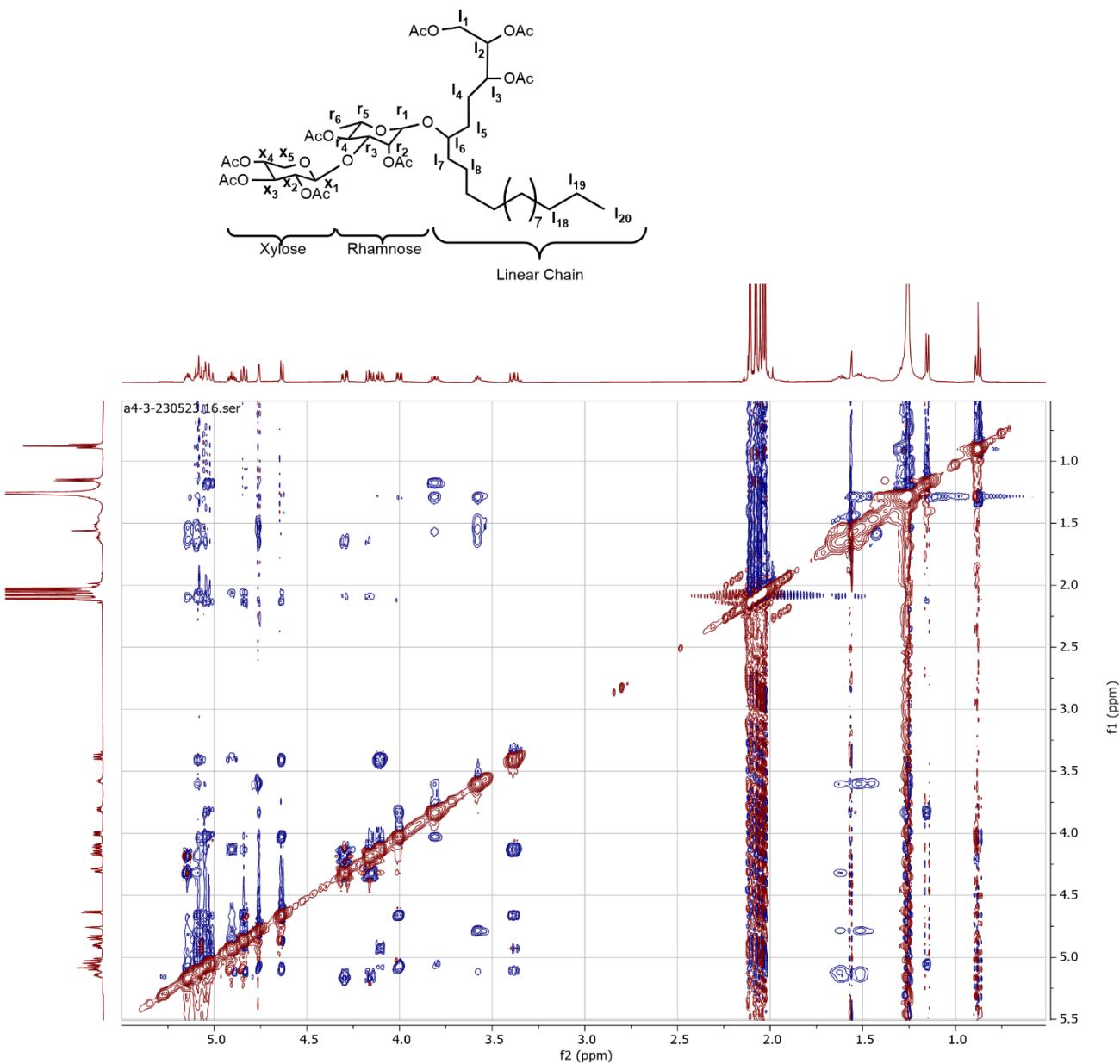
(b) 10-80 ppm range.



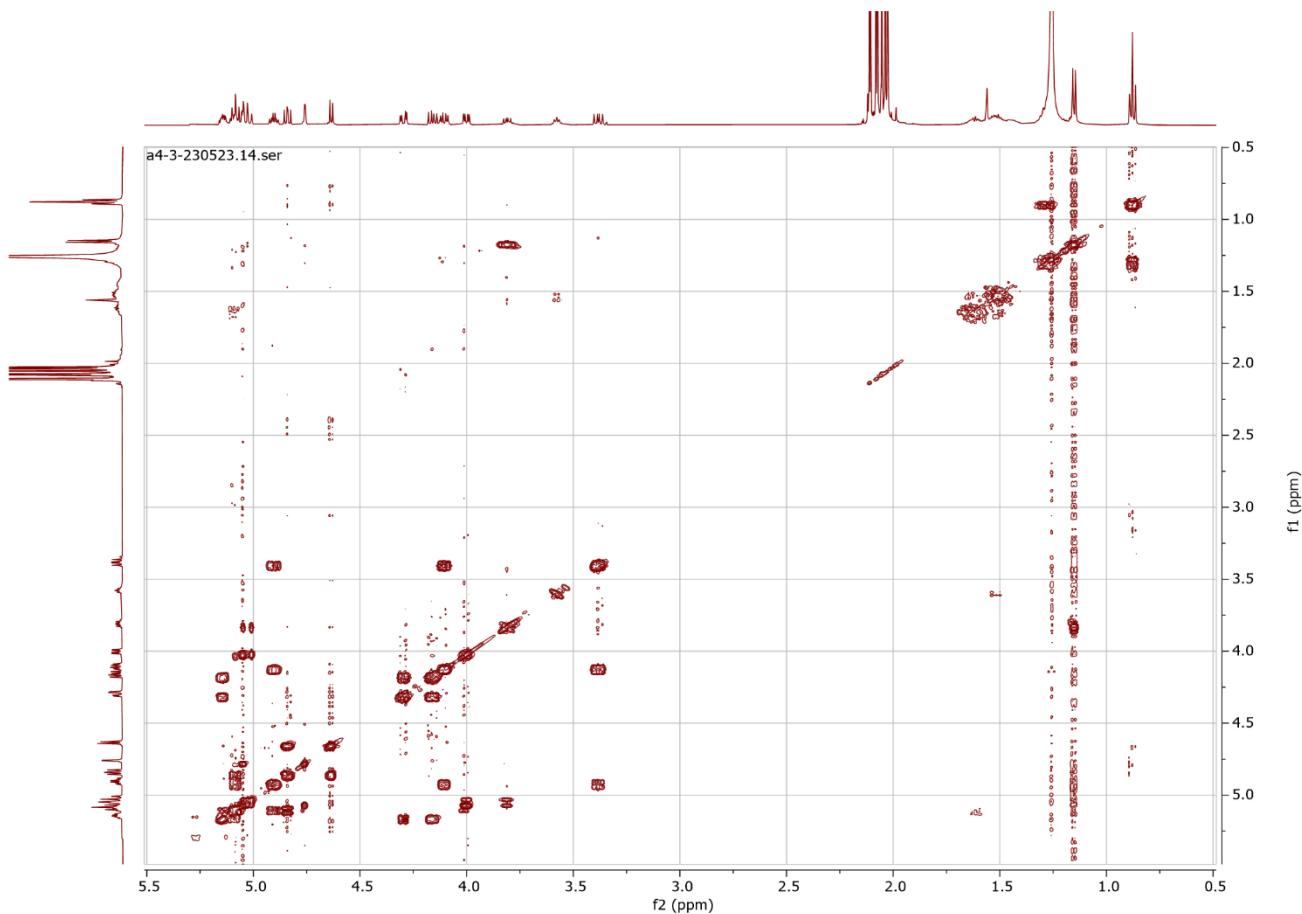
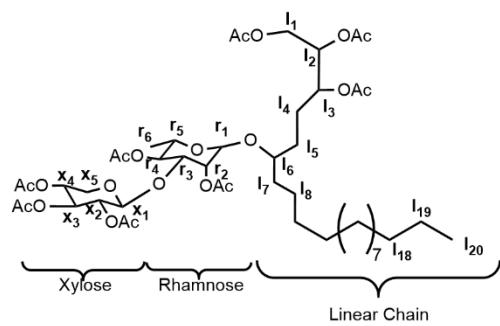
Supplementary Fig 8: ¹H-¹³C one bond (¹J) correlation pattern (HSQC, 500 MHz, CDCl₃) for compound **II** (¹H: 0.5-5.5 ppm range/ ¹³C: 10-110 ppm range).



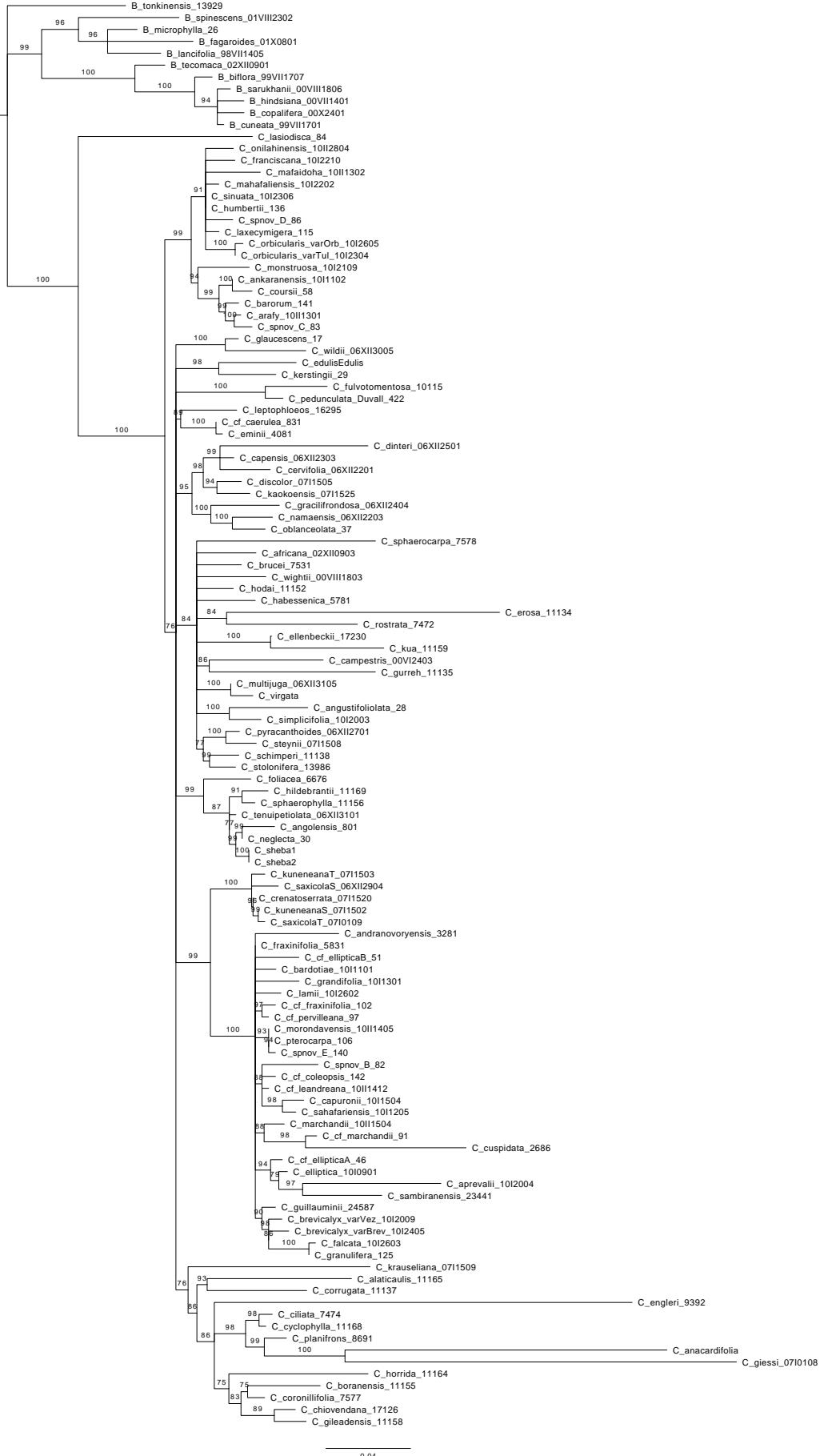
Supplementary Fig 9: ^1H - ^{13}C long range ($^{2,3}J$) correlation pattern (HMBC, 500 MHz, CDCl_3) for compound **II**. ^1H : 0.5-5.5 ppm range/ ^{13}C : 10-180 ppm range



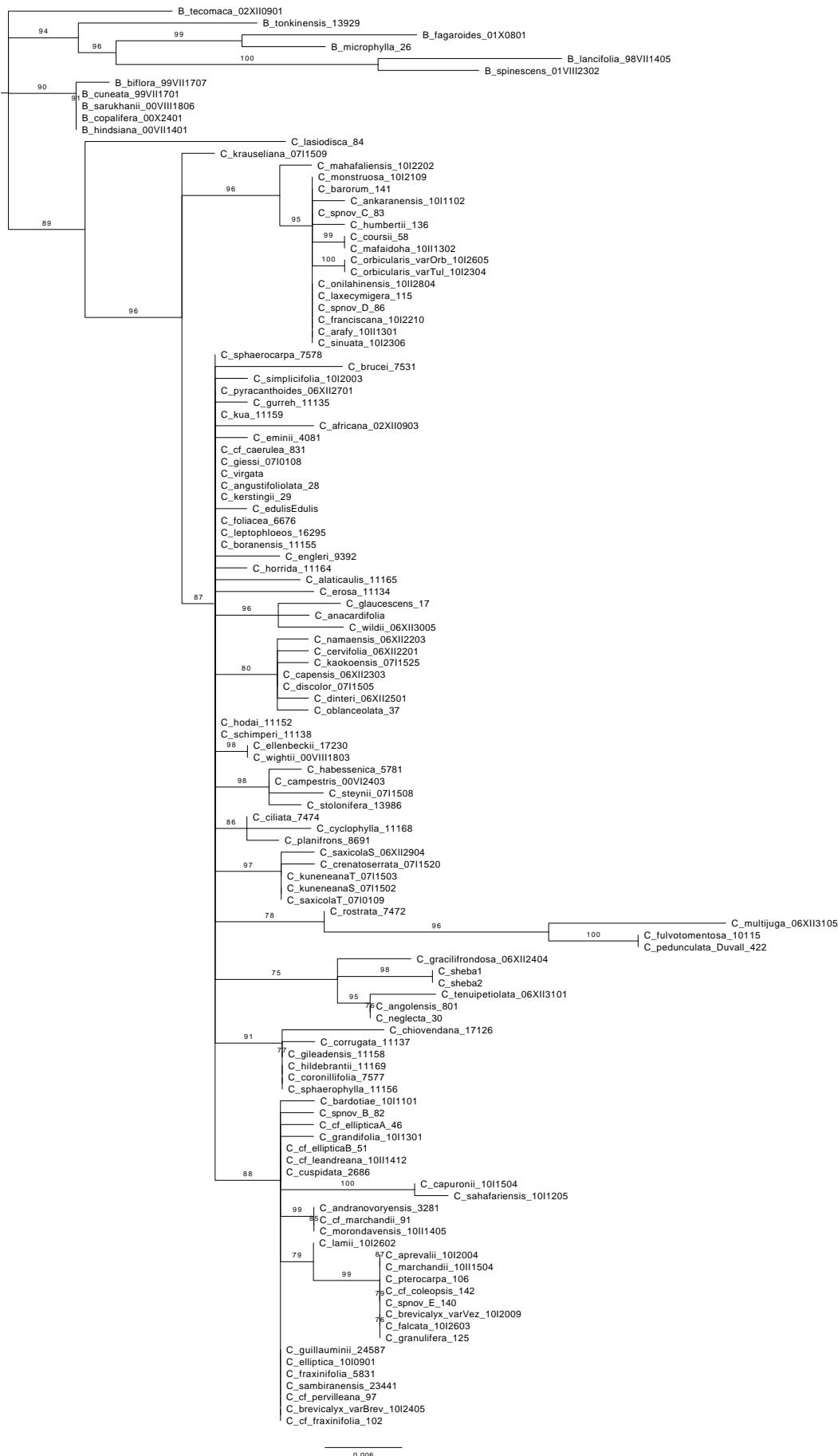
Supplementary Fig 10: ^1H - ^1H NOESY correlation pattern (500 MHz, CDCl_3) for compound **II** (0.5-5.5 ppm range).



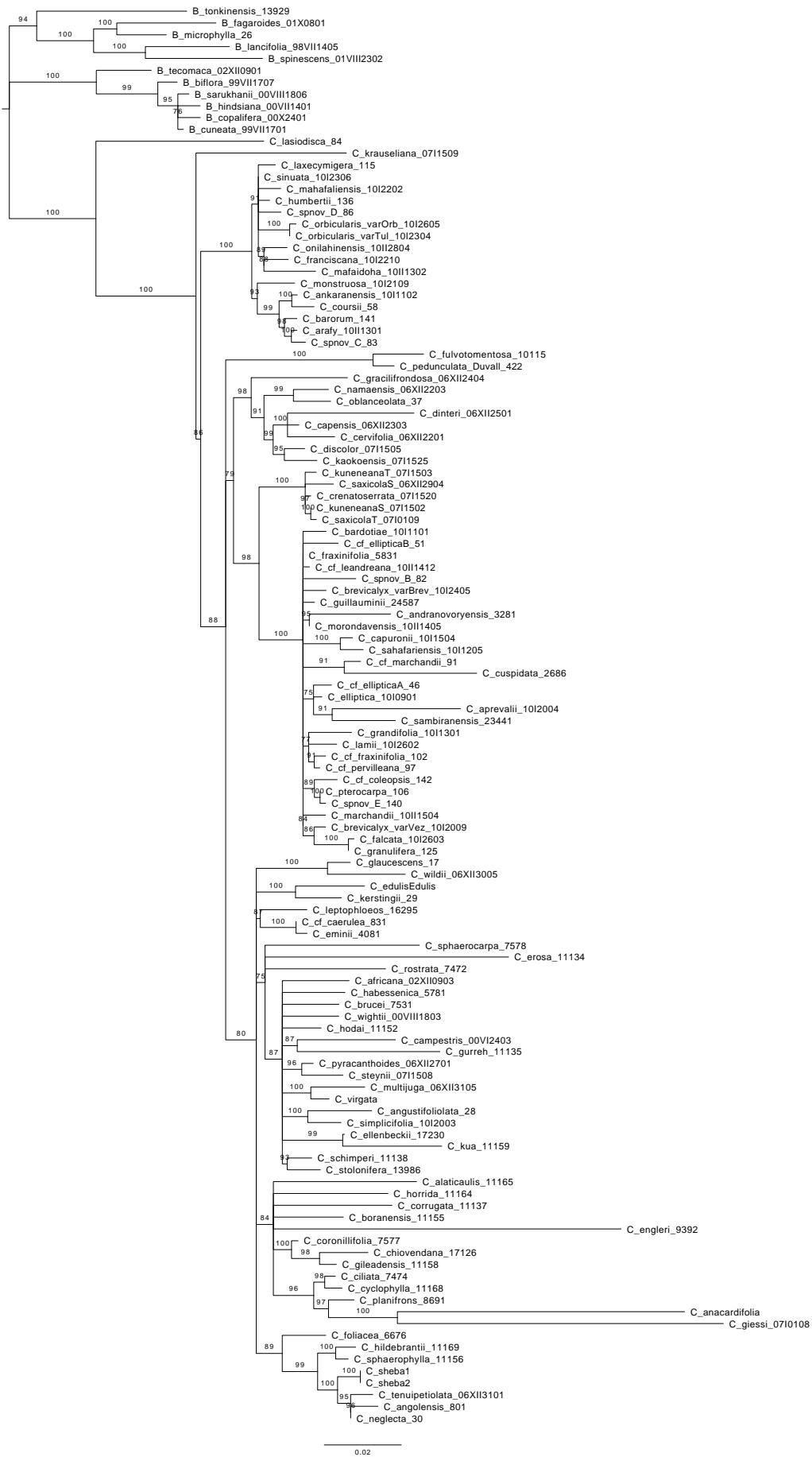
Supplementary Fig 11: ^1H - ^1H COSY correlation pattern (500 MHz, CDCl_3) for compound **II** (0.5–5.5 ppm range).



Supplementary Fig 12: Phylogeny of *Commiphora* based on maximum likelihood analysis of nuclear DNA sequence data. Tree topology is the 75% majority rule consensus phylogram of 1000 ultrafast bootstrap replicate trees.



Supplementary Fig 13: Phylogeny of *Commiphora* based on maximum likelihood analysis of chloroplast DNA sequence data. Tree topology is the 75% majority rule consensus phylogram of 1000 ultrafast bootstrap replicate trees.



Supplementary Fig 14: Phylogeny of *Commiphora* based on maximum likelihood analysis of combined nuclear and chloroplast DNA sequence data. Tree topology is the 75% majority rule consensus phylogram of 1000 ultrafast bootstrap replicate trees.

Supplementary References

- 1.**Bronk Ramsey, C. Development of the radiocarbon calibration program. Radiocarbon 43, 355-363 (2001) DOI: <https://doi.org/10.1017/S0033822200038212>
- 2.**Bronk Ramsey, C.Bayesian analysis of radiocarbon dates. Radiocarbon 51, 337-360 (2009)
DOI: <https://doi.org/10.1017/S0033822200033865>
- 3.** Reimer, P. J. et al. The IntCal20 Northern Hemisphere radiocarbon age calibration curve (0–55cal kBP). Radiocarbon. 62, 4, 725–757 (2020). DOI:[10.1017/RDC.2020.41](https://doi.org/10.1017/RDC.2020.41)