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Supporting Information

Total Synthesis of (±) Commiphoranes C-D and their Epimers

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1. General Information

¹H NMR and ¹³C NMR were recorded on Mercury 400, Bruker AV500, AV600, JEOL ECZ400S spectrometer. Coupling constants are given in Hz and chemical shifts are expressed as δ values in ppm. The following multiplicity abbreviations are used: (s) singlet, (d) doublet, (t) triplet, (q) quartet, (m) multiplet. ESI-HRMS data were measured on Thermo Exactive Orbitrap plus spectrometer. All the chemicals were purchased from commercial sources: Sigma-Aldrich Chemical Co., Arcos Chemical Co., and J&K Chemical Co. with the purity of more than 95%. Solvents were dried according to standard procedures when needed. Flash column chromatography was performed on Biotage Isolera one or carried out on silica gel (200–300 mesh). IR spectra were recorded on a Thermo Nicolet 5700 FT-IR microscope Centaurµs spectrophotometer.



2. ¹H, ¹³C, NOE and 2D-NMR Spectral Copies

¹³C NMR spectra of compound 8



 ^1H NMR spectra of compound (±) 5



¹³C NMR spectra of compound (±) 5



¹³C NMR spectra of compound (±) 4



¹³C NMR spectra of compound (±) 10







¹³C NMR spectra of compound (±) 12



¹³C NMR spectra of compound (±) 14



¹³C NMR spectra of compound (±) 3



¹³C NMR spectra of compound (±) **15**



¹³C NMR spectra of compound (±) 16



¹³C NMR spectra of compound (±) 2



¹³C NMR spectra of compound **1a**



3.812 3.793 3.773 3.506

2. 375 2. 281 2. 101 2. 101 2. 101 2. 101 2. 071 3. 067 3.

.071

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pdata/1 Bruker AVANCEIII 400 20210427 & & & PROTON2 CDC13 D:\\ DATA-2021 22

 $< \frac{5,831}{5,805}$

¹³C NMR spectra of compound **1b**







H-H cosy spectra of compound **1c**







¹H NMR spectra of compound **1d**



¹³C NMR spectra of compound **1d**







NOE spectra of compound 1d

	natural commiphorane C		1a	
position	$\delta_{\rm H}(600{\rm M})$	$\delta_{\rm C}(150{\rm M})$	$\delta_{ m H}$ (400M)	$\delta_{\rm C}(100{\rm M})$
1		121.2		121.0
2		53.2		53.1
3	4.24 (t-like, 4.8)	71.1	4.24 (t, 5.1)	70.9
4	a 2.20 (dd, 1.4, 13.1),	39.8	a 2.19 (dd, 1.6, 12.0),	39.6
	b 2.10 (dd, 4.8, 13.1)		b 2.09 (dd, 4.8, 12.6)	
5	2.18 (m)	32.3	2.18 (m)	32.1
6	5.91 (d, 10.0)	74.7	5.91 (d, 9.7)	74.5
7		138.6		138.4
8	6.81 (s)	119.1	6.81 (s)	119.0
9		140.0		139.8
10	6.86 (s)	110.5	6.86 (s)	110.4
11		153.5		153.4
12		179.7		179.5
13	1.53 (s)	19.6	1.53 (s)	19.5
14	1.03 (d, 6.4)	19.6	1.03 (d, 6.2)	19.4
15	2.38 (s)	22.3	2.38 (s)	22.1
-OAc		170.1		170.0
-OAc	2.29 (s)	21.0	2.29 (s)	20.8
-OH				

Table 1 The comparison of $^1\mathrm{H}$ and $^{13}\mathrm{C}$ NMR data of compound 1a and reported natural commiphorane C

	natural commiphorane D		1b	
position	$\delta_{\rm H}(600{\rm M})$	$\delta_{\rm C}(150{\rm M})$	$\delta_{\rm H}$ (400M)	$\delta_{\rm C}(100{\rm M})$
1		122.2		122.2
2		49.9		49.9
3	3.79 (t, 7.4)	69.9	3.73-3.85 (m)	69.9
4	2.08 (dd, 7.4, 9.9,	40.2	2.03-2.14 (m)	40.3
	2H)			
5	1.87 (m)	37.7	1.79-1.97 (m)	37.7
6	5.82 (d, 10.5)	74.5	5.82 (d, 10.4)	74.5
7		139.0		139.0
8	6.81 (s)	119.5	6.81 (s)	119.5
9		140.1		140.2
10	6.88 (s)	110.6	6.88 (s)	110.6
11		152.8		152.8
12		181.5		181.5
13	1.58 (s)	14.6	1.58 (s)	14.6
14	1.06 (d, 6.7)	19.8	1.06 (d, 6.7)	19.8
15	2.38 (s)	22.2	2.37 (s)	22.2
-OAc		170.1		170.1
-OAc	2.28 (s)	21.0	2.28 (s)	21.0
-OH	3.51(s)		3.51(s)	

Table 2: the comparison of $^1\mathrm{H}$ and $^{13}\mathrm{C}$ NMR data of compound 1b and reported natural commiphorane D

Compounds 1a-1d and (±) 11 inhibited the fibrogenic action of TGF- β 1 in rat renal proximal tubular cells.

Cell lysates were immunoblotted with antibodies against collagen I as indicated. After marker, the compounds were **1b**, **1c**, **1d**, **1a** and (±) **11** (25 μ M), and **1b**, **1c**, **1d**, **1a** and (±) **11** (50 μ M), DMSO with TGF- β , DMSO without TGF- β , GW: GW788388 in turn.

Cell lysates were immunoblotted with antibodies against α -SMA, and α -tubulin as indicated. After marker, the compounds were **1b**, **1c**, **1d**, **1a** and (±) **11** (25 μ M), and **1b**, **1c**, **1d**, **1a** and (±) **11** (50 μ M), DMSO with TGF- β , DMSO without TGF- β , GW: GW788388 in turn.