## A New Sesquiterpene Lactones Glucoside from Notoseris psilolepis

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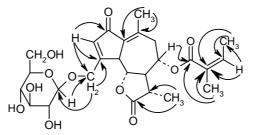
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**Abstract:** A new sesquiterpene lactone glucoside, notoserolide C (1), was isolated from the whole plant of *Notoseris psilolepis* Shih. By means of spectral analysis including MS, NMR (<sup>1</sup>H NMR, <sup>13</sup>C NMR, <sup>1</sup>H-<sup>1</sup>H COSY, HMQC, HMBC) and X-ray diffraction, the structure of notoserolide C was established as cichorioside B angelate.

**Keywords:** *Notoseris psilolepis*, asteraceae, notoserolide C.

As a part of an investigation on plant of the genus *Notoseris* (Asteraceae), an endemic genus of seed plants of China<sup>1</sup>, we studied the chemical constituents of *N. psololepis* Shih. A new sesquiterpene lactone glucoside, named notoserolide C (1) was isolated from the methanolic extract of the whole plant of *N. psilolepis* by repeated column chromatography (MCI GEL, normal and reversed phase silica gel). This compound showed antibacterial activity against *Bacillu cereus*.

Figure 1. The key HMBC correlations for 1



Notoserolide C (1), was obtained from methanol as colorless crystals, mp 156~8°C (dec.), [  $\alpha$  ] $_D^{25}$  -55° (c 0.07, MeOH). Its molecular formula was assigned as  $C_{26}H_{34}O_{11}$  by HRFABMS ([M-H]  $^{-}$  m/z 521.1992, Calcd. 521.1964). The IR spectrum of 1 showed the presence of hydroxyl (3413 cm $^{-1}$ ) and carbonyl (1765, 1717, 1686 cm $^{-1}$ ) groups. On acid hydrolysis, glucose was detected by TLC. Since the signal of the anomeric proton of the glucose appeared at  $^{\delta}_{\rm H}$  4.40 with coupling constant 8 Hz, this sugar moiety should be  $^{\beta}$ -orientated. Besides the glucosyl moiety, the  $^{1}\text{H-NMR}$  spectrum exhibited a methyl

doublet at  $\,^{\delta}$  1.19 (3H, d, J = 7 Hz), a vinyl methyl signal at  $\,^{\delta}$  2.44 (3H, s) , an olefinic proton signal at  $\,^{\delta}$  6.61 (1H, br s) and angeloyl signals at  $\,^{\delta}$  6.23 (1H, q, J = 7 Hz), 2.01 (3H, d, J = 7 Hz) and 1.92 (3H, s). The  $^{13}$ C-NMR spectrum was similar to that of cichorioside  $\,^{2}$  except for five signals due to the angeloyl moiety $^{3}$ . The long-range heteronuclear correlations (**Figure 1**) between H-15 and the anomeric carbon of the glucose, between H-8 and carbonyl carbon of the ester moiety revealed that the glucose moiety is located at C-15 and angeloyl at C-8. Analysis of  $^{1}$ H- $^{1}$ H COSY, HMQC and HMBC spectra allowed proton and carbon signals of  $\,^{1}$  to be assigned as **Table 1**. Accordingly,  $\,^{1}$  can be represented as cichoroside B angelate. The structure was unambiguously confirmed by X-ray analysis $^{3}$ .

 $\delta_{C}$  $\delta_{\,\underline{c}}$  $\delta_{\,\underline{H}}$  $\delta_{\,\underline{\,}H}$ C C 1 134.0(s) ester moiety 2 197.1(s) 16 167.8(s) 134.9(d) 128.5(s) 3 6.61, br s 17 171.7(s) 18 140.8(d) 6.23, q, 7Hz 5 19 2.01, d, 7Hz 3.87, d, 8Hz 49.7(d) 16.1(q) 6 82.2(d) 3.84, m 20 1.92, s 20.7(q)7 59.4(d) 2.55, m sugar moiety 8 71.5(d) 5.00, m 1' 104.1(d) 4.40, d. 8Hz 2' 9 45.6(t)2.90, m 75.1(d) 10 148.6(s) 3' 78.0(d) 11 41.7(d) 2.67, m 4 71.5(d) 5 78.0(d) 179.1(s) 12 13 15.2(q) 1.19, d, 7Hz 62.7(t)3.67, dd, 11, 7Hz 3.88, d, 11Hz 14 21.5(q) 2.44, s 4.80, 4.87, d, 17Hz 15 69.6(t)

**Table 1**.  $^{1}$ H-(500 MHz) and  $^{13}$ C-(125 MHz) NMR spectral data for **1** (CD<sub>3</sub>OD)

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## **References and Notes**

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- 3. Crystallographic parameter of 1 have been deposited in the editorial office of CCL.

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