## Saluenolide A, A Novel Eremophilanolide from Senecio saluenensis

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**Abstract:** From *Senecio saluenensis*, a novel eremophilanolide Saluenolide A was isolated. Its structure was elucidated by 2D-NMR technique and X-ray diffraction.

Keywords: Sesquiterpene, eremophilanolide, Senecio, X-ray, cytotoxicity.

The crude petroleum ether extract of *S. saluenensis* showed 78% inhibition to KB cell on  $100 \mu \text{mol/L}^1$ . Nine known compounds and a new eremophilanolide, saluenolide A (1) was isolated from this extract.

Saluenolide A (1) showed 22 carbon resonances, combined by DEPT data and the molecular ion peak exhibited at m/z 340, the molecular formula of 1 could be deduced as  $C_{22}H_{30}O_8$ . Three typical methyl signals at  $\delta$  1.92 (br s), 1.24 (s) and 0.95 (d, J=6.8 Hz) suggested that this compound is an eremophilanolide<sup>2,3</sup>. Considering the presence of three esteric carboxylic resonances at  $\delta$  170.9, 170.3 and 165.4, two ester moieties exist in the molecule. Scrutiny on the  $^1H$  and  $^{13}C$ -NMR spectra disclosed the existence of an acetoxy group as well as a senecioyl in the molecule. Two hydroxyls appeared at  $\delta$  5.09 and 3.86 could be exchanged by  $D_2O$ . One olefinic proton appeared at  $\delta$  5.67 (br s) could be assigned as H-2 in the senecioyl moiety. Two oxygen-bearing methines exhibited at  $\delta$  5.72 (br s, H-3) and 4.85 (d, J=2.4 Hz, H-6). Furthermore, the ketal carbon resonance at  $\delta$  103.1 indicated the presence of the 8-hydroxy-8, 12-olide moiety<sup>2</sup>.

Another hydroxyl should locate at C-10, since this oxygen-bearing carbon signal appeared at  $\delta$  74.2. The HMBC as well as the NOESY experiment revealed the presence of a 3 $\beta$ -acetoxy group and a 6 $\beta$ -senecioyl moiety. The absolute stereochemistry of **1** was finally disclosed by X-ray diffraction experiment utilized a Mo atom initiation. **1** was evaluated for its inhibitory activity against the growth of HL-60, A-549 and KB cell lines *in vitro* with IC<sub>50</sub> values of  $5.6 \times 10^{-5}$ ,  $4.8 \times 10^{-5}$  and  $6.1 \times 10^{-4}$  mol/L, respectively.

Structuee of Saluenolide A

X-ray structure of Saluenolide A

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