고추(*Capsicum annuum* L.) 잎으로부터 Acyclic Diterpenoids의 분리 경희대학교: 조진경, 이재웅<sup>1</sup>, 이대영, 이도경, 류하나, 이윤형<sup>1</sup>, 백남인\*

## Acyclic Diterpenoids from the Leaves of Capsicum annuum L.

Graduate School of Biotechnology & Plant metabolism Research Center, Kyung Hee University Department of Life Science, <sup>1</sup>Department of Horicultural Biochemistry Kyung Hee university

Jin-Gyeong Cho, Jae-Woong Lee<sup>1</sup>, Dae-Young Lee, Do-Gyeong Lee, Ha-Na Lyu, Youn-Hyung Lee<sup>1</sup>, Nam-In Baek\*

## **Objectives**

Capsicum species are very important plants used worldwide as vegetable foods, spices, and external medicine. A number of studies have been carried out on *C. annuum* fruit, but there are few studies on leaves. *C. annuum* leaves were very familiar to Korean as seasoned vegetable in spring and summer season. Though *C. annuum* leaves were reported as inhibitory effect on anti-complementary activity, anti-mutagenic activity, anti-microbial activity, DPPH scavenging activity, and anti-tyrosinase activity, any phytochemical investigation has not been reported so far. In this study, the authors reported the isolation and identification of three terpenoids from *C. annuum* leaves.

## Materials and Methods

The IR spectrum was obtained with a Perkin Elmer Spectrum One FT-IR spectrometer,  $CaF_2$  window in MeOH (Buckinghamshire, England). EI-MS was recorded on a JEOL JMSAX-505-WA.  $^1\text{H-NMR}$  (400 MHz) and  $^{13}\text{C-NMR}$  (100 MHz) spectra were recorded on a Varian Unity Inova AS-400 FT-NMR spectrometer.

The leaves of C annuum. were extracted with 80% aqueous MeOH, and the concentrated extract was partitioned with EtOAc, n-BuOH and H<sub>2</sub>O, successively. From the EtOAc fraction, three acyclic diterpenoids were isolated through the repeated silica gel and ODS column chromatographies.

## Results

From the EtOAc fraction, three acyclic diterpenoids were isolated through the repeated silica gel and ODS column chromatographies. According to the results of physico-chemical data including NMR, MS and IR, the chemical structures of the compounds were determined as *a*-tocopherol (1), phytol (2), and so on.

Corresponding author: Nam-In Baek E-mail: nibaek@khu.ac.kr Tel: 031-201-2661