

A New Compound from *Dalbergia odorifera* T. Chen

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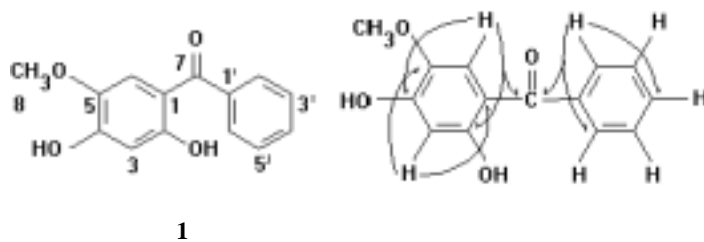
Abstract: A new compound, 2, 4-dihydroxy-5-methoxybenzophenone was isolated from the root and heartwood of *Dalbergia odorifera* T. Chen.

Keywords: Benzophenone, *Dalbergia odorifera*.

The plant *Dalbergia odorifera* T. Chen (Leguminosae) is a high tree, the root and heartwood of this plant is a Chinese traditional medicine named Jiangxiang. It is indigenous to Guangdong province.

Compound **1** was isolated from the EtOAc extract of plant *Dalbergia odorifera* T. Chen as yellow crystalline powder, mp: 191.2-192.5 °C. EI-MS (m/z): 244 (96), 243 (92), 167 (100), 166 (65), 123 (23), 105 (55), 77 (95). IR ν (KBr) cm^{-1} : 3300 (hydroxy), 1637 (β -unsaturated carbonyl), 1597, 1571 and 1512 (aromatic ring). Its spectrum showed UV λ_{max} nm: 252, 287 (aromatic ring), 362. $^1\text{H-NMR}$ (400MHz) (CDCl_3 , δ): 3.96 (3H, s, H-8), 5.24 (1H, s, -OH), 6.54 (1H, s, H-3), 7.10 (1H, s, H-6), 7.47 (2H, ddd, $J=7.6, 7.3, 2.2\text{Hz}$, H-3', 5'), 7.45 (1H, dd, $J=7.3, 2.2\text{Hz}$, H-4'), 7.63 (2H, dd, $J=7.6, 2.2\text{Hz}$, H-2', 6'), 12.48 (1H, s, -OH). The $^{13}\text{C-NMR}$ (100MHz) (CDCl_3) and DEPT spectra showed signal at δ 56.2 (C-8), δ 99.9 (C-3), δ 111.8 (C-1), δ 116.7 (C-6), δ 128.3 (C-2', 6'), δ 128.7 (C-3', 5'), δ 131.4 (C-4'), δ 137.8 (C-4), δ 138.3 (C-1'), δ 153.6 (C-5), δ 160.1 (C-2), δ 200.0 (C-7). The partial structure was elucidated from the HMBC spectrum; most of which are shown by arrows in **Figure 2**. In addition, the following $^{13}\text{C-}^1\text{H}$ long-range correlation was also observed between H-3 and C-1, C-5; H-6 and C-2, C-4, C-7; H-8 and C-5; H-2' and C-4', C-6', C-7 signals. The $^1\text{H-}^1\text{H}$ COSY and HMQC spectra sported assignment of all proton signals. Thus, the structure was determined as 2, 4-dihydroxy-5-methoxy-benzo-phenone.

Figure 1 Structure of compound **1** **Figure 2** The HMBC correlation of compound **1**



References

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