

Phytochemical communication

Constituents of *Caesalpinia pyramidalis*

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Abstract

Two new glycosyl phenylpropenoid acids, 4-*O*- β -glucopyranosyloxy-(*Z*)-7-hydroxycinnamic acid (**1**) and 4-*O*- β -glucopyranosyloxy-(*Z*)-8-hydroxycinnamic acid (**2**), besides lupeol and aghatisflavone, were isolated from the leaves of *Caesalpinia pyramidalis*. © 2000 Elsevier Science B.V. All rights reserved.

Keywords: *Caesalpinia pyramidalis*; Phenylpropenoid glycosides; Biflavones; Triterpenes

Plant. *Caesalpinia pyramidalis* Tul. (Caesalpinaceae), dried leaves, collected in Valente (BA), Brazil by Prof. W.A. Lopes, in August 1993, and identified by Prof. L. Scardino (IB-UFBA). Voucher specimen is deposited under number 0240291 at Herbário Alexandre Leal Costa of Instituto de Biologia of Universidade Federal da Bahia.

Uses in traditional medicine. For fever and stomach diseases, and as diuretic [1].

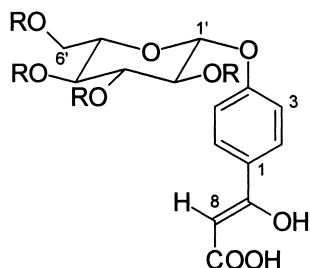
Previously isolated constituents. No report.

New-isolated constituents. 4-*O*- β -Glucopyranosyloxy-(*Z*)-7-hydroxycinnamic acid (**1**, yield 0.001%), 4-*O*- β -glucopyranosyloxy-(*Z*)-8-hydroxycinnamic acid (**2**) [2]

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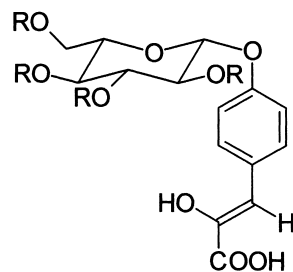
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(0.00073), isolated as acetyl derivatives **1a** and **2a**, respectively; lupeol [3] (0.034) and aghatisflavone [4] (0.076).



1 R=H

1a R=COCH₃



2 R=H

2a R=COCH₃

4-O-β-(2',3',4',6'-Tetraacetyl)glucopyranosyloxy-(Z)-7-hydroxycinnamic acid (1a). UV max (MeOH): 218 (ϵ 61963), 274 (39840); (+ AlCl₃): 222 (61868), 282 (36409), 330 (20767); (+ AlCl₃ + HCl): 224 (60033), 282 (35331), 330 (20818) nm; IR bands (KBr): 3415, 2953, 1754, 1658, 1228, 846 cm⁻¹; ¹H-NMR (300 MHz, CDCl₃): δ 7.79 (2H, *d*, *J* 8.5 Hz, H-2 and H-6), 7.04 (2H, *d*, *J* 8.5 Hz, H-3 and H-5), 6.57 (1H, *s*, H-8), 6.39 (1H, *d*, *J* 8.7 Hz, H-1'), 5.68–4.02 (4H, *m*, H-2'-H-5'), 4.38 (1H, *dd*, *J*₁ 12.5, *J*₂ 2.8, H-6'), 4.22 (1H, *dd*, *J*₁ 12.5, *J*₂ 2.0, H-6'), 2.11 (6H, *s*, CH₃CO), 2.04 (6H, *s*, CH₃CO), 12.81 (1H, *s*, OH); ¹³C-NMR (75 MHz, CDCl₃): 123.41 (C-1), 128.14 (C-2 and C-6), 116.43 (C-3 and C-5), 162.69 (C-4), 159.44 (C-7), 98.71 (C-8), 182.08 (C-9), 104.33 (C-1'), 67.78, 70.73, 73.66 and 73.67 (C-2'-C-5'), 61.23 (C-6'), 19.87, 20.40, 20.48 and 20.50 (4 × CH₃CO), 170.43, 170.29, 169.21 and 168.40 (4 × COO).

4-O-β-(2',3',4',6'-Tetraacetyl)glucopyranosyloxy-(Z)-8-hydroxycinnamic acid (2a). UV max (MeOH): 226 (ϵ 42794), 270 (27380); (+ AlCl₃): 234 (87890), 280 (41487), 304 (44211); (+ AlCl₃ + HCl): 234 (88656), 280 (42324), 304 (45116) nm; IR bands (KBr): 3419, 1751, 1654, 1240, 840 cm⁻¹; ¹H-NMR (300 MHz, CDCl₃): δ 7.94 (2H, *d*, *J* 8.4 Hz, H-2 and H-6), 7.37 (2H, *d*, *J* 8.4 Hz, H-3 and H-5), 6.65 (1H, *s*, H-7), 6.41 (1H, *d*, *J* 8.7 Hz, H-1'), 5.69–4.02 (4H, *m*, H-2'-H-5'), 4.39 (1H, *dd*, *J*₁ 12.5, *J*₂ 1.5, H-6'), 4.22 (1H, *dd*, *J*₁ 12.5, *J*₂ 1.5, H-6'), 2.12 (6H, *s*, CH₃CO), 2.04 (6H, *s*, CH₃CO), 12.81 (1H, *s*, OH); ¹³C-NMR (75 MHz, CDCl₃): 128.69 (C-1), 127.48 (C-2

and C-6), 122.74 (C-3 and C-5), 162.71 (C-4), 127.30 (C-7), 153.60 (C-8), 181.95 (C-9), 105.94 (C-1'), 67.74, 70.68, 73.61 and 73.62 (C-2'-C-5'), 61.26 (C-6'), 20.96, 20.44, 20.38 and 19.91 ($4 \times \underline{\text{CH}}_3\text{CO}$), 170.37, 170.11, 169.13 and 168.67 ($4 \times \underline{\text{COO}}$).

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References

- [1] Inventário de Plantas Medicinais do Estado da Bahia. Governo do Estado da Bahia (SEPLANTEC), 1979:334.
- [2] Shigemori H, Sakai N, Miyoshi E, Shizuki Y, Yamamura S. *Tetrahedron* 1990;46:383.
- [3] Mahato SB, Kundu AP. *Phytochemistry* 1994;37:1517.
- [4] Chari VM, Ilyas M, Wagner H et al. *Phytochemistry* 1977;16:1273.