*Phytochemistry* Vol 28, No. 6, pp. 1787–1789, 1989.
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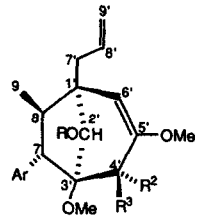
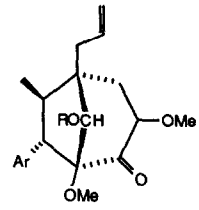
# ERRATUM

C. D. Dodson *et al.* (1987) *Phytochemistry* **26**, 2037. The publishers regret the structures in the above paper were incorrectly drawn, with the bridges toward the oxygen rather than the carbon. The correct structures are given below.




1a-4a




5a-7c

# CORRIGENDA

F. Balza *et al.* (1988) *Phytochemistry* **27**, 1421. The authors regret that in the Experimental, the $\delta$c (ppm) valves for $CDCl_3$ and $Me_2CO$-$d_6$ were reversed; it should read $Me_2CO$-$d_6$ ($\delta c = 29.8$ ppm) and $CDCl_3$ ($\delta c = 77.0$ ppm).

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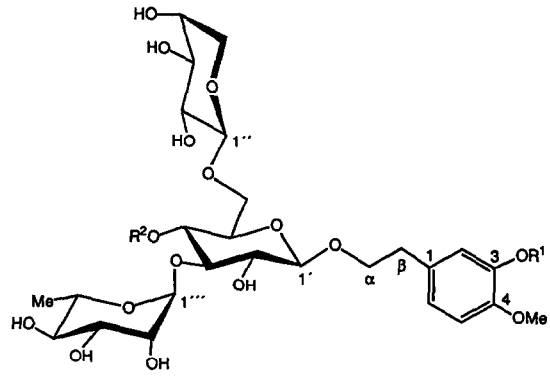
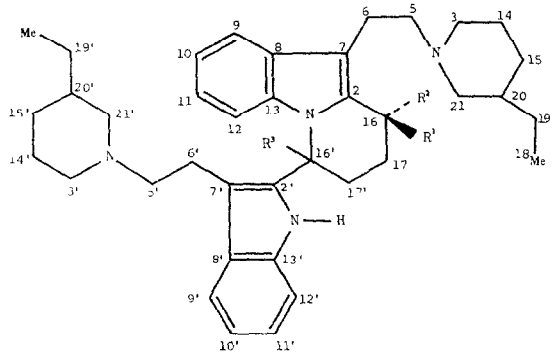
I. Calis *et al.* (1988) *Phytochemistry* **27**, 1465. The authors regret a mistake in the arabinose part of the structure formula, the correct structure is given below.




| | R¹ | R² |
|---|---|---|
| 1 | H | caffeoyl |
| 2 | H | feruloyl |
| 3 | H | H |
| 4 | Me | H |

A.-ur-Rahman and K Laman (1988) *Phytochemistry* **27,** 1926. The authors regret a mistake in the structure of compounds **1** and **2** The correct structures are as follows



**1** $R^1 = R^3 = H$, $R^2 = CO_2Me$

**2** $R^1 = CO_2Me$, $R^2 = R^3 = H$

Lines 10–12 in column 2, p. 1927 should now read, "in **1** the $CO_2Me$ was α-oriented while in **2** the $CO_2Me$ was β-oriented" Line 5 should read, "C-16 from the β-isomer to the α-isomer" and line 14 should read, "the $CO_2Me$ group being presented at C-16′"

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R Tanaka and S. Matsunaga (1988) *Phytochemistry* **27,** 2274. The authors regret a mistype in the $^{13}C$ NMR chemical shift value of the C-2 carbon signal for compound **7a** listed in Table 3. It should be δ27 41 not δ24.41

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S Yu *et al* (1988) *Phytochemistry* **27,** 2887 The authors regret that 6-hydroxypulchellin-4-*O*-angelate was reported as a new compound from *Gaillardia pulchella*; in fact this compound has been previously reported from the same species by Professor S. Inayama and co-workers. S Inayama *et al* (1983) *Heterocycles* **20,** 1501

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N Gopalsamy *et al* (1988) *Phytochemistry* **27,** 3593 In the formula shown on p. 3593, the stereochemistry of the methyl group at C-14 should be α and not β, otherwise the aglycone depicted is not tormentic acid

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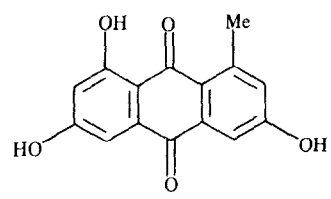
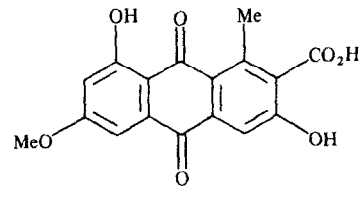
A. A Ali *et al* (1989) *Phytochemistry* **28,** 281. Incorrect formulae were given in the above paper, the correct structures are given below.

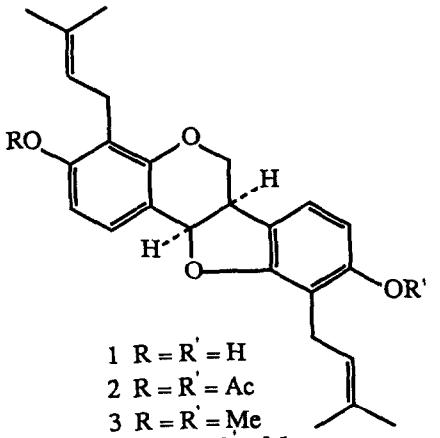


**1**

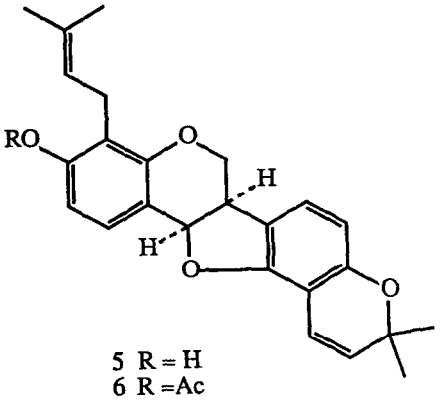


**2**

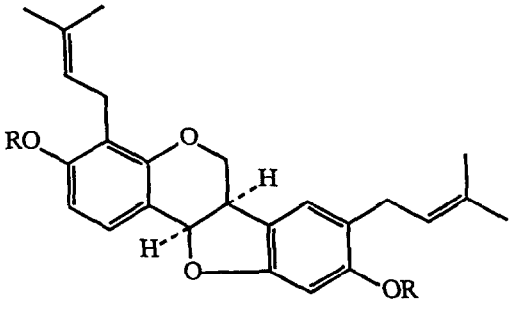
L. A. Mitscher *et al.* (1988) *Phytochemistry* **27**, 3449. The following structures were omitted from the above paper.



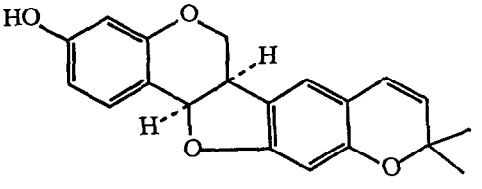
1 R = R' = H
2 R = R' = Ac
3 R = R' = Me
4 R = H, R' = Me



5 R = H
6 R = Ac



7 R = H
8 R = Ac



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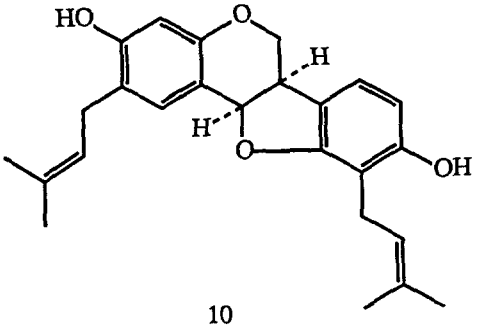

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