

TWO NEW STEREOISOMERIC COUMARINS FROM THE LEAF EXTRACTIVES  
OF *CALOPHYLLUM CORDATO-OBLONGUM*  
THW. (GUTTIFERAE)

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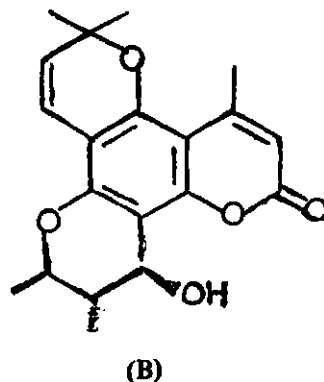
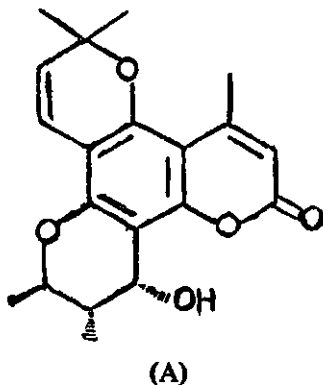
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As a part of continuing studies of the extractives of the plant family Guttiferae, the foliar constituents of *Calophyllum cordato-oblongum* Thw. were investigated. The light petroleum extract of the leaves had four coumarins, two of which have been characterised. Both had molecular formula  $C_{20}H_{22}O_6$ . Both had similar mass spectral fragmentations having the base peak at  $m/z$  327( $M^+ - CH_3$ ). This fragmentation is typical of the compounds having 2, 2-dimethyl-2H pyran system(1) (2) The presence of the following protons were confirmed from

the  $^1H$  n.m.r. data:  $-CH=CH(J=10Hz)$ ;  $CH_3-C=CH$ ;  $-CH-OH$ ;  $2xCH_3-CH$ ;  $-\overset{O}{\parallel}C=CH=C$ . The i.r. absorption at  $1735\text{ cm}^{-1}$  ( $\gamma C=O$ ) for both compounds suggested that they are coumarins. This proposal is quite reasonable since the u.v. spectra of both compounds were identical and were very similar to those reported for a number

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of 5, 7-dioxygenated coumarins. Besides, both of these compounds gave the same oxidised product. The spectroscopic and chemical evidence show that these coumarins are 11, 12-dihydro-12 $\alpha$ -hydroxy-4, 6,6,10 $\beta$ , 11 $\alpha$ -pentamethyl-2H, 6H, 10H-benzo [1, 2, 6; 3, 4, 6'; 5, 6, 6'']-tripyran-2-one (A) and 11,12-dihydro-12 $\beta$ -hydroxy-4, 6, 6, 10 $\beta$ , 11 $\alpha$  -pentamethyl-2H, 6H, 10H-benzo [1, 2, 6; 3, 4, 6'; 5, 6, 6'] -tripyran-2-one (B).



### References

1. Barnes, C. S., *et al Tetrahedron Letters*, 281 (1963).
2. Stout, G. E. and Stevens, K. L., *J. Org. Chem.*, 29, 3604 (1964).