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## 3,3'-Benzylidenebis(4-hydroxy-6-methylcoumarin). Erratum

An error in the paper by Vijayalakshmi *et al.* [Acta Cryst. (2002), E**58**, o659–o660] is corrected. The last sentence of page o659 should read "The C21–C22 distance of 1.534 (2) Å is longer than an unstrained  $Csp^3-C_{ar}$  bond, but is in the range characteristic of sterically crowded structures ... ".

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