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

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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 . . . ".