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1-Acetyl-3-ethyl-*r*-2,*c*-6-di-2-furylpiperidin-4-one. Corrigendum

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In the paper by Balamurugan, Thiruvalluvar, Manimekalai, Selvaraju & Maruthavanan [*Acta Cryst.* (2006), E**62**, o2005–o2006], the two first sentences in the *Abstract* contain errors. The correct text is "In the title molecule, $C_{17}H_{19}NO_4$, the piperidine ring adopts a chair conformation. The acetyl group in the 1-position and the ethyl group in the 3-position have equatorial and axial orientations, respectively".