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Addition/Correction

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Correction to "A Cyclochiral Conformational Motif Constructed Using a Robust Hydrogen-Bonding Network"

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Page 13645. In Figure 3b, the unit of $\Delta S^{\ddagger}$ was incorrectly given as "$\Delta S^{\ddagger} = -10.8$ kcal/mol". It should be "$\Delta S^{\ddagger} = -10.8$ cal/mol·K". The corrected version of Figure 3 is presented. The authors apologize for this error.

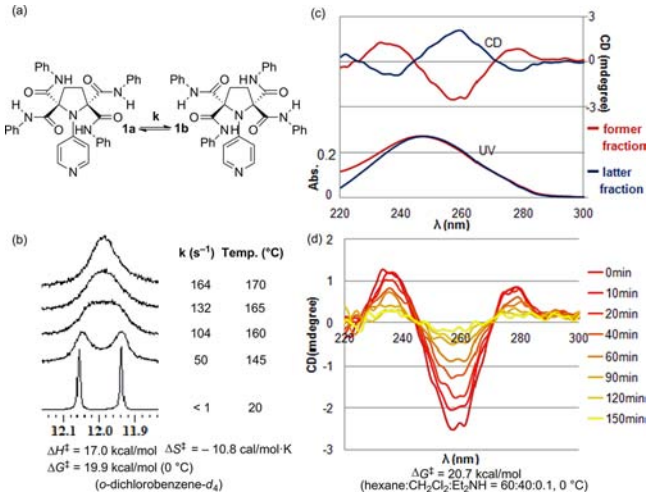

Figure 3. Analysis of the racemization barrier of **1**. (a) Interconversion of the enantiomers of **1** through rotation of the amide moieties. (b) Coalescence of the amide protons observed in a VT NMR study in *o*-dichlorobenzene-d_4. (c) CD spectra of both enantiomers in hexane:CH_2Cl_2:Et_2NH (60:40:0.1). (d) A trace of racemization was observed based on the CD spectra.

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