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* Author to whom correspondence should be addressed

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# CORRIGENDA

In the paper by K. C. Nicolaou et al. published in *Chem. Eur. J.* **2001**, *7*, 5359 – 5371, the enantiomeric structures of the Diels–Alder components in the transition states $\mathbf{TS_a}$ and $\mathbf{TS_b}$ are required to account for the observed enantiomers **8**, (−)-**5**, **8′**, and **5′** shown in Scheme 9.

In the paper by J. Jiang, N. Kobayashi, and D. K. P. Ng et al. in *Chem. Eur. J.* **2001**, *7*, 5059 – 5069, the true space group for all the double-deckers $\mathbf{4}\cdot 0.5\,C_6H_{12}$, **6**, $\mathbf{9}\cdot 0.5\,C_6H_{12}$, and **14** is *Pnam* (No. 62; interchange of the *b* and *c* axes as reported in the paper leads to the standard symbol *Pnma*). The $[M^{III}(nc)(oep)]$ molecule lies on a crystallographic mirror plane that contains the lanthanide atom, two *trans*-related isoindole N donor atoms of the nc ring, and a pair of *trans*-related *meso* CH groups on the oep ligand. In the complexes $\mathbf{4}\cdot 0.5\,C_6H_{12}$ and $\mathbf{9}\cdot 0.5\,C_6H_{12}$, the chair-shaped cyclohexane molecule has half-site occupancy with two of $CH_2$ groups lying in a mirror plane. The authors apologize for these errors.

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