

Corrigenda

The First Internally Functionalized Chiral [2.2]Metacyclophanes

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The last line in the abstract should be replaced by:

'enantiomers of (**2c–h**) are compared with those of the less strained parent skeleton (**2b**).'

Crystal data for (**2d**), 5th to 7th line should be as follows:

' cm^{-1} , 3946 reflections with $|F| > 4\sigma(F)$ were used for the structure solution (direct methods, Enraf-Nonius SDP¹³) and refinement (253 parameters, Enraf-Nonius SDP¹³), non-hydrogen atoms were refined' . . .

The last two lines in the text to Table 1 should be replaced by:

'hydrogens H_i and inner methyl groups of the [2.2]phanes (**2b**) to (**2i**) [X = NTos (Tos = *p*-SO₂C₆H₄Me)] and (**2b**).'

The text to Figure 3 should be as follows:

Figure 3: CD spectra of (**2c–h**), compared to (**2b**). (**2c**): $\Delta\epsilon_{227} = +60.8$, $\Delta\epsilon_{276} = -4.7$; (**2d**): $\Delta\epsilon_{227} = +80.4$, $\Delta\epsilon_{276} = -6.5$, $\Delta\epsilon_{291} = -4.0$; (**2e**): $\Delta\epsilon_{232} = +31.2$, $\Delta\epsilon_{278} = -7.7$; (**2f**): $\Delta\epsilon_{232} = +23.5$, $\Delta\epsilon_{270} = -4.4$, $\Delta\epsilon_{295} = -3.2$; (**2g**): $\Delta\epsilon_{232} = +23.4$, $\Delta\epsilon_{276} = -3.2$; (**2h**): $\Delta\epsilon_{235} = +11.5$, $\Delta\epsilon_{265} = -5.9$, $\Delta\epsilon_{295} = -1.8$.

The first lines in the 3rd paragraph on p. 1759 should be replaced by:

'The separation or enrichment of the enantiomers of the new chiral compounds (**2c–h**) was successfully accomplished using' . . .

Ref. 5(b) should be replaced by:

'R. H. Mitchell, in "Cyclophanes," eds. P. M. Keehn and S. M. Rosenfeld, Academic Press, New York, 1983;' . . .