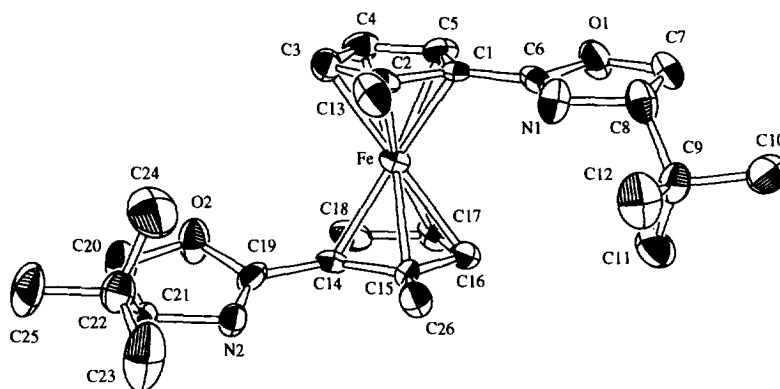


## CORRIGENDUM

S0957-4166(96)00130-9

Wanbin Zhang, Yoichi Adachi, Toshikazu Hirao and Isao Ikeda, Highly diastereoselective *ortho*-lithiation of 1,1'-bis(oxazoliny)ferrocene directed to  $C_2$ -symmetric chiral ligands, *Tetrahedron: Asymmetry*, **1996**, 7, 451–460, S0957-4166(96)00028-6

On page 453, Figure 1, the crystal structure of **6b** was enantiomerically misprinted and the corrected Figure 1 should be as follows:



**Figure 1** Crystal structure of **6b** (ORTEP, ellipsoids at the 30% probability level). Selected bond length [Å] and bond and torsion angle [°]: C(1)-C(2) 1.42 (1), C(2)-C(13) 1.50 (2), C(1)-C(6) 1.51 (1), O(1)-C(6) 1.39 (1), N(1)-C(6) 1.24 (1), O(1)-C(7) 1.45 (2), N(1)-C(8) 1.48 (1), C(7)-C(8) 1.52 (2), C(8)-C(9) 1.53 (2); C(1)-C(2)-C(13) 128 (1), C(2)-C(1)-C(6) 123 (1), O(1)-C(6)-N(1) 118 (1), O(1)-C(6)-C(1) 110 (1), N(1)-C(6)-C(1) 130 (1), C(5)-C(1)-C(6) 126 (1); N(1)-C(6)-C(1)-C(2) 5(1), C(1)-C(6)-O(1)-C(7) 179.3(9), C(1)-C(6)-N(1)-C(8) 178 (1), C(6)-N(1)-C(8)-C(9) -122 (1), C(6)-C(1)-C(2)-C(13) 4 (1).

On page 453, paragraph 4 line 2, '(S)-planar' should be replaced by "(R)-planar".