

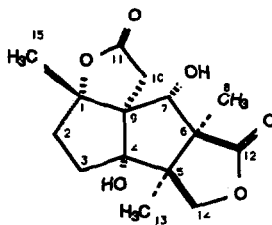
### CORRIGENDA

R. Pellicciari, B. Natalini and M. Marinozzi, D-3,4-'Cyclopropylglutamate' isomers as NMDA receptor ligands: synthesis and enantioselective activity, *Tetrahedron Lett.*, 1990, 31, 139

The absolute configurations of the *trans*-isomers D-CGA A and D-CGA B have been inadvertently inverted. According to the correct structures depicted on page 139, they should read as follows:  
on page 141, line 11 from the top, D-CGA B, 8, (2*R*,3*R*,4*R*)  
instead of (2*R*,3*S*,4*S*)  
on page 141, line 15 from the top, D-CGA A, 7, (2*R*,3*S*,4*S*)  
instead of (2*R*,3*R*,4*R*)

I. Kouno, K. Mori, N. Kawano and S. Sato, Structure of anisactone A; a new skeletal type of sesquiterpene from the pericarps of *Illicium anisatum*, *Tetrahedron Lett.*, 1989, 30, 7451

The authors wish to apologize for the critical error in the configuration at C-7 of anisactone A. It should be 7 $\alpha$ -OH and not 7 $\beta$ -OH:



(1) anisactone A