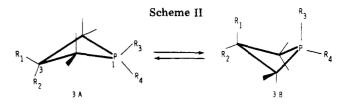
S. Alunni, E. Baciocchi,* P. Perucci, and R. Ruzziconi: Influence of the Steric Requirements of the Nucleophile on the Transition State Structure of E2 Reactions. A Kinetic Study of the Elimination from 1-Bromo-2-arylethanes and 1-chloro-1-phenyl-2-arylethanes Promoted by Sodium 2,6-Di-*tert*-butylphenoxide.

Page 2415. Table I, line 1 should read as follows: nucleophile, M (0.023); k_2 , M^{-1} s⁻¹ (6.22×10^{-4}) . A. Fitzgerald,* J. A. Campbell,* G. D. Smith, C. N. Caughlan,

A. Fitzgerald, * J. A. Campbell, * G. D. Smith, C. N. Caughlan, and S. E. Cremer: Solid-State Studies on Crowded Molecules. Crystal and Molecular Structures of 2,2,3-Trimethyl-1-phenylphosphetane 1-Oxide and 2,2,3,3,4-Pentamethyl-1-phenylphosphetane 1-Oxide.

Page 3517. First column. The figure shown below (Scheme II in the text) is needed for an understanding of the last paragraph in the first column of page 3517.



Mitsuru Imuta and Herman Ziffer*: Preparation and Absolute Stereochemistry of Isomeric Pyridylethanols and *threo*-Di(2-pyridyl)ethanediol.

Page 3532. First column. Structure 4 in Scheme II is incorrect; the portion shown as "C-CH₃" should be "C-COOCH₃", i.e., 4 is (S,S)-(+)-dimethyl diacetyltartrate, as designated correctly in the text.