

Corrigenda**Di- and Tetra-nuclear Complexes with Bis(diphenylphosphino)amide and Bis(diphenylphosphino)methanide as Bi- and Tri-dentate Ligands. X-Ray Structures of $[(\text{Ph}_3\text{P})(\text{O}_3\text{ClO})\text{AgN}(\text{Ph}_2\text{PAuPPh}_2)_2\text{NAg}(\text{OCIO}_3)(\text{PPh}_3)]$ and $[(\text{C}_6\text{F}_5)\text{AuCH}(\text{Ph}_2\text{PAuPPh}_2)_2\text{CAu}(\text{C}_6\text{F}_5)]$**

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J. Chem. Soc., Chem. Commun., 1986, 509.

On p. 510 the footnote 'Crystal data' contains an error: the space group of compound (4) is given as $P1$ but should be $P\bar{1}$.

The following information is missing from this article: the atomic co-ordinates for this work are available on request from the Director of the Cambridge Crystallographic Data Centre, University Chemical Laboratory, Lensfield Rd., Cambridge CB2 1EW. Any request should be accompanied by the full literature citation for this communication.

Selective Formation of Substituted Cyclopentane Derivatives from Hexa-1,5-dienes via Oxidative Cyclization in the Presence of $\text{Pd}(\text{OAc})_2\text{-MnO}_2\text{-Benzoquinone}$ as Catalyst

Thomas Antonsson, Andreas Heumann, and Christina Moberg

J. Chem. Soc., Chem. Commun., 1986, 518.

On page 519, the structure of compound (16) should be as shown below.

