The structure of **5b** was solved by a direct method (MULTAN 82). The final refinements were accomplished by a full-matrix least-squares method with anisotropic thermal parameters for non-hydrogen atoms. All calculations were performed on a MICRO-VAX II computer by using the TEXSAN program system.

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Supporting Information Available: Tables of data collection and processing parameters, positional and thermal parameters, bond lengths, and bond angles for **5b** (4 pages). Ordering information is given on any current masthead page.

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Additions and Corrections

1997, Volume 16

Wolfgang Ahlers, Bodo Temme, Gerhard Erker,* Roland Fröhlich, and Frank Zippel: Formation, Structure, and Dynamic Behavior of a Novel Dinuclear Cationic μ-2,4-Hexadiyne Bis(zirconocene) Complex.

Page 1443. Reference 19 is incorrect. The correct literature citation is as follows.

(19) Strauss, D. A.; Zhang, C.; Tilley, T. D. *J. Organomet. Chem.* **1989**, *369*, C13.

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