

Erratum: “Assessment of conventional density functional schemes for computing the polarizabilities and hyperpolarizabilities of conjugated oligomers: An ab initio investigation of polyacetylene chains” [J. Chem. Phys. 109, 10489 (1998)]

Benoît Champagne, Eric A. Perpète, Stan J. A. van Gisbergen, Evert-Jan Baerends, Jaap G. Snijders, Chirine Soubra-Ghaoui, Kathleen A. Robins, and Bernard Kirtman

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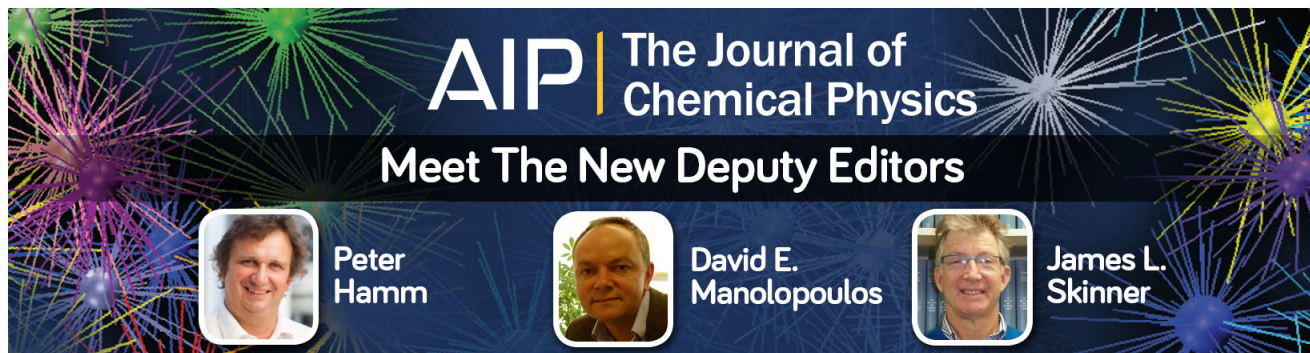
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ERRATA

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Benoît Champagne and Eric A. Perpète

*Laboratoire de Chimie Théorique Appliquée, Facultés Universitaires Notre-Dame de la Paix,
rue de Bruxelles 61, B-5000 Namur, Belgium*

Stan J. A. van Gisbergen and Evert-Jan Baerends

*Section Theoretical Chemistry, Vrije Universiteit, De Boelelaan 1083, 1081 HV Amsterdam,
The Netherlands*

Jaap G. Snijders

*Department of Chemical Physics, University of Groningen, Nijenborgh 4, 9747 AG Groningen,
The Netherlands*

Chirine Soubra-Ghaoui and Kathleen A. Robins

Department of Chemistry, University of Nevada, Las Vegas, Nevada 89154

Bernard Kirtman

Department of Chemistry, University of California, Santa Barbara, California 93106

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The external electric field used to obtain Fig. 2 was 20×10^{-4} a.u. rather than 20×10^4 a.u.