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)]

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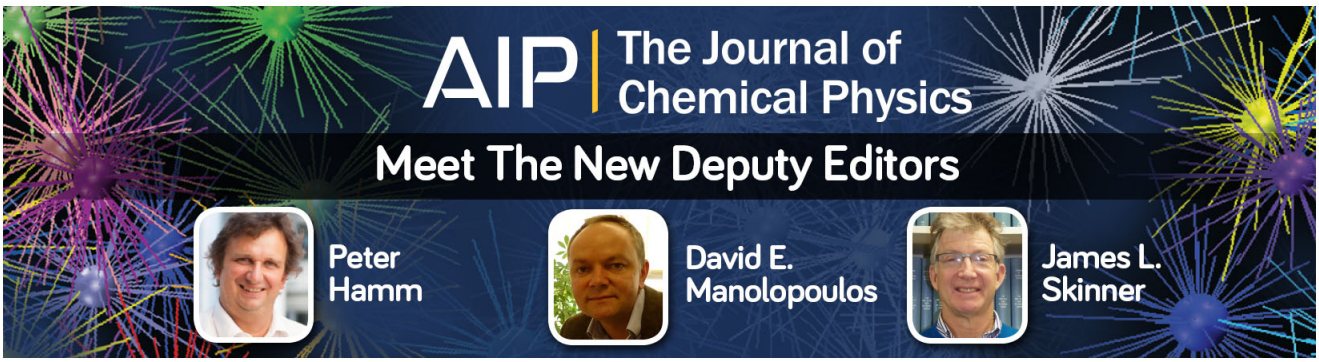
Citation: The Journal of Chemical Physics **110**, 11664 (1999); doi: 10.1063/1.479106
View online: http://dx.doi.org/10.1063/1.479106
View Table of Contents: http://scitation.aip.org/content/aip/journal/jcp/110/23?ver=pdfcov
Published by the AIP Publishing



LETTERS TO THE EDITOR

The Letters to the Editor section is divided into three categories entitled Notes, Comments, and Errata. Letters to the Editor are limited to one and three-fourths journal pages as described in the Announcement in the 1 January 1999 issue.

ERRATA

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[S0021-9606(99)00623-6]

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

