Additions and Corrections

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J. D. Smith and T. P. Hanusa*: Trends in the Structures and Energetics of the Group 14 Metallocenes $(C_5H_5)_2M$ (M = Si−Pb): A Density Functional Theory Study

Page 3059. The footnotes to Table 2 are incorrect and should read as follows.

[a]Energies in hartrees; energies in parentheses are kcal mol^{-1} equivalents. The lowest absolute energies (au) for each metallocene are as follows: Si, −390.631 332 (ECP); Si, −676.201 697 (all-electron); Ge, −681.469 692; Sn, −601.202 045: Pb, −579.780 213. [b]With pseudopotential basis set on Si. [c]With all-electron basis set (aug-cc-pVTZ) on Si.

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