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*Sarom S. Leang, Federico Zahariev, and Mark S. Gordon*

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*Jun Li, Yimin Wang, Bin Jiang, Jianyi Ma, Richard Dawes, Daiqian Xie, Joel M. Bowman, and Hua Guo*

[J. Chem. Phys. 136, 041103 \(2012\)](#)

### **Perspective on density functional theory**

*Kieron Burke*

[J. Chem. Phys. 136, 150901 \(2012\)](#)

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*V. Holten, C. E. Bertrand, M. A. Anisimov, and J. V. Sengers*

[J. Chem. Phys. 136, 094507 \(2012\)](#)

### **Accurate ab initio potential energy surface, thermochemistry, and dynamics of the Cl(<sup>2</sup>P, <sup>2</sup>P<sub>3/2</sub>) + CH<sub>4</sub> → HCl + CH<sub>3</sub> and H + CH<sub>3</sub>Cl reactions**

*Gábor Czako and Joel M. Bowman*

[J. Chem. Phys. 136, 044307 \(2012\)](#)

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