

Long-range exchange interactions in DFT and their significance in chemical reactions

Thursday, 22 June 2017 15:15 (1h 15m)

The significance of long-range exchange interactions is presented from the viewpoint of DFT in quantum chemistry [1] in the first part, and then, it is shown focusing on chemical reactions in the second part. So far, we have developed the long-range corrected (LC) DFT [2] and have applied it to a wide variety of chemical and physical properties [3]. As a result, we have confirmed that the long-range exchange interactions are required to calculate various types of the properties: e.g. charge transfer excitations [4], van der Waals bonds [5], nonlinear optical properties [1] and so forth. Orbital energies may be the most significant property that LC-DFT makes it possible to calculate quantitatively [6]. Since orbital energies are the solution of the Kohn-Sham equation, this indicates that the long-range correction essentially improves DFT (or exactly the generalized DFT). In the first part, I will briefly review our past studies on LC-DFT.

Recently, we are investigating chemical reactions using the quantitative orbital energies. As the exact orbital energies are proven to inhere, LC-DFT orbital energies hardly vary dependent on occupation numbers. Chemical reactions usually proceed through charge transfers in the initial processes. We found that LC-DFT orbital energies are kept almost constant in the initial processes of many reactions, and then, they rapidly increase toward the products [7]. We have recently developed an orbital energy-based reaction theory as the modification of conceptual DFT [8]. I will present this topic in the second part.

1. T. Tsuneda, "Density Functional Theory in Quantum Chemistry" (Springer, 2014).
2. H. Iikura, T. Tsuneda, T. Yanai, and K. Hirao, *J. Chem. Phys.*, 115, 3540 - 3544, 2001.
3. T. Tsuneda and K. Hirao, *WIREs Computational Molecular Science* 4, 375 - 390, 2014.
4. T. Tsuneda and K. Hirao, "Time-Dependent Density Functional Theory", in *Theoretical and Quantum Chemistry at the 21st Century Dawn End* (Apple Academic Press, 2017).
5. T. Tsuneda and T. Taketsugu, " π -Stacking on Density Functional Theory: A review", in *π -Stacked Polymers and Molecules*, Ed. T. Nakano, 245 - 270 (Springer, 2013).
6. T. Tsuneda, J.-W. Song, S. Suzuki, and K. Hirao, *J. Chem. Phys.* 133, 174101(1 - 9), 2010.
7. T. Tsuneda and R. K. Singh, *J. Comput. Chem.* 35, 1093 - 1100, 2014.
8. T. Tsuneda, "Chemical reaction analyses based on orbitals and orbital energies", *Int. J. Quantum Chem. Special Issue on Theoretical Chemistry in Japan* 115, 270 - 282, 2015.

Presenter: Prof. TSUNEDA, Takao (University of Yamanashi)