FRISTERIS Hub Meeting **Vol.43** Vol.27 発表者: Nguyen Tuan Hung Assist. Prof. (東北大学 学際科学フロンティア研究所/物質材料・エネルギー)

Density-functional theory: A powerful tool for interdisciplinary materials science

*Language: English

Summary

Density-functional theory (DFT) is a computational method used in materials science, chemistry, and condensed matter physics to study the electronic structure of materials. It is based on quantum mechanics and can predict accurate results, which can compare directly with experimental data. Thus, DFT has become an important tool for predicting the properties of materials, such as their electronic structure, chemical reactivity, crystal lattices, and spectroscopic properties. This talk will introduce how to use the DFT in interdisciplinary materials science. We also show the hands-on tutorial for Quantum ESPRESSO, a wide-used open software for DFT calculation, on personal computers.

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2023.



11:00 -12:00

ハイブリッド開催(学際科学フロンティア研究所セミナー室 & Online)

