19th International Workshop on Computational Physics and Material Science: Total Energy and Force Methods | (smr 3266)

Contribution ID : 6 Type : not specified

Electronic structure of transition metal compounds by variational cluster approximation

Wednesday, 9 January 2019 11:50 (0:30)

Content

Summary

 $\textbf{Presenter(s)}: \quad \text{EDER, Robert (Karlsruhe Institute of Technology, Wuerttemberg, Germany)}$

Session Classification: Session 2 - Strong Correlations in Quantum Materials