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John M. Wills, Mebarek Alouani, Per Andersson, Anna Delin, Olle Eriksson, Oleksiy Grechnyev

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## Full-Potential Electronic Structure Method: 167 (Springer **Series in Solid-State Sciences)**

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Full-Potential Electronic Structure Method: 167 (Springer Series in Solid-State Sciences) John M. Wills, Mebarek Alouani, Per Andersson, Anna Delin, Olle Eriksson, Oleksiy Grechnyev This book covers the theory of electronic structure of materials, with special emphasis on the usage of linear muffin-tin orbitals. Methodological aspects are given in detail as are examples of the method when applied to various materials. Different exchange and correlation functionals are described and how they are implemented within the basis of linear muffin-tin orbitals. Functionals covered are the local spin density approximation, generalised gradient approximation, self-interaction correction and dynamical mean field theory.



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