

Ab initio Calculations of Spin-Wave Excitation Spectra from Time-Dependent Density-Functional Theory

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Schriften des Forschungszentrums Jülich
Reihe Schlüsseltechnologien / Key Technologies

Band / Volume 38

ISSN 1866-1807

ISBN 978-3-89336-786-3

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Schlüsseltechnologien / Key Technologies
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