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Eleventh International Workshop on Computational Physics and Material Science: Total Energy and Force Methods | (smr 1491)

Friday 17 January 2003

<u>Ab-initio Simulations of Molecular Fluids: Structural and Electronic Properties and the Influence of a Magnetic Field</u> - Main Building Main Lecture Hall (09:00-09:30)

time	title	presenter
09:00	Ab-initio Simulations of Molecular Fluids: Structural and Electronic Properties and the Influence of a Magnetic Field	GIULIA GALLI