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Conference on Single Molecule Magnets and Hybrid Magnetic Nanostructures | (smr 1664)

Friday 01 July 2005

Density-Functional-Based Simulation of Molecular Magnets: Anisotropy and Exchange Hamiltonians - Main Building Main Lecture Hall (09:00-09:45)

time	title	presenter
09:00	Density-Functional-Based Simulation of Molecular Magnets: Anisotropy and Exchange Hamiltonians	MARK PEDERSON