

Poster Session

Symposium on „Functional Transition Metal Compounds & Multiferroics“
Cologne 26.09.05 – 28.09.05

A'AVO(PO₄)₂: The first examples of frustrated ferromagnetic square lattice systems (A', A = Pb, Ba, Sr, Zn)

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Crystal structure and magnetic properties of La_{2-x}Sr_xCoO₄

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Role of the structural distortions in the phase diagrams of RE_{1-x}A_xTiO₃ (A=Sr, Ca)

H. Roth, A. Komarek, M. Cwik, W. D. Stein, N. Schnittner, A. El-Filali, M. Kriener, T. Zabel, A. Freimuth, M. Braden, T. Lorenz

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Multiferroic Crystals: Crystal Growth and Characterization

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Halides of the outer and inner transition metals / Complex oxides of the

3d transition metals

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Growing thin films with control of oxygen stoichiometry

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Controlling orbital occupation and spin orientation in transition metal oxide thin films

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Pressure-induced metal-insulator transition in $RNiO_3$

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Transport Properties of low-dimensional Cuprates, Vanadates, and Nickelates

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Thermoelectric Properties of doped $LaCoO_3$

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Thermodynamic Properties and Metamagnetism of Ca_{2-x}Sr_xRuO₄

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Thermal Properties of Multiferroic Materials

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Optical spectroscopy on highly correlated transition metal compounds I

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Optical spectroscopy on highly correlated transition metal compounds II

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Magnetism in 214-Ruthenates

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Magnetic excitations in single-layer manganates

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Unusual spin state and spin-state transitions in cobaltates

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Phase transitions and orbital occupation in partially filled t_{2g} systems

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Resonant soft x-ray diffraction

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Spectroscopy from stripe order in $La_{1.8}Sr_{0.2}NiO_4$

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L Dynamical Correlations in Ferromagnetic Transition Metal Compounds: a view from LDA+DMFT

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Low dimensional quantum spin systems with ring exchange and spin-phonon coupling

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Almost integrable models

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Exact ground states of quantum spin systems in $d>1$ dimensions

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Disorder and quantum fluctuation in effective field theories of highly correlated materials

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Field Tuned Quantum Critical Points in Antiferromagnetic Metals

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Chiral Helices in a Metal: MnSi

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Ab initio and Cluster Model Calculations for Strongly Correlated Oxide (1)

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Ab initio and Cluster Model Calculations for Strongly Correlated Oxide (2)

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