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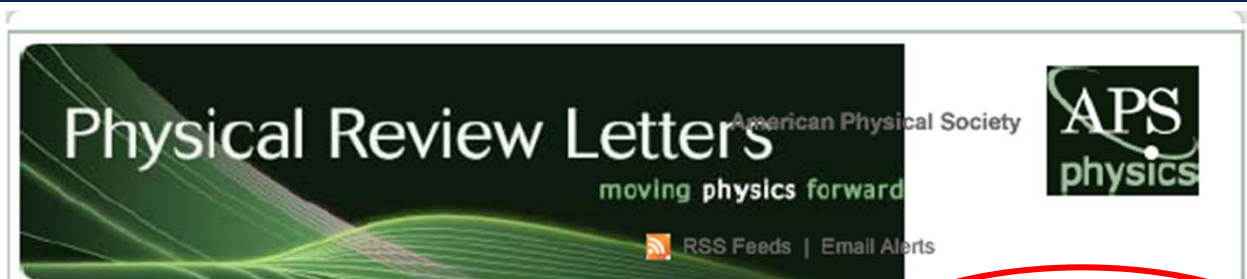
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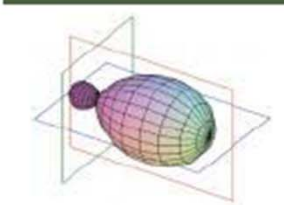
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[J.-S. Zhou](#) and [J. B. Goodenough](#)

Texas Materials Institute, 1 University Station, ETC 9.104, University of Texas at Austin, Austin, Texas 78712, USA

(Received 13 April 2006; published 19 June 2006)

We report the refinement of x-ray powder diffraction together with magnetic and thermal conductivity measurements made on the entire family of $RMnO_3$ perovskites prepared by melt growth or under high pressure. Analysis of the data has identified the origin of the transition from type-A to type-E magnetic order as a competition between t -orbital and e -orbital spin-spin interactions within each Mn-O-Mn bond in the (001) planes, the e -orbital interactions decreasing with decreasing R^{3+} -ion size.

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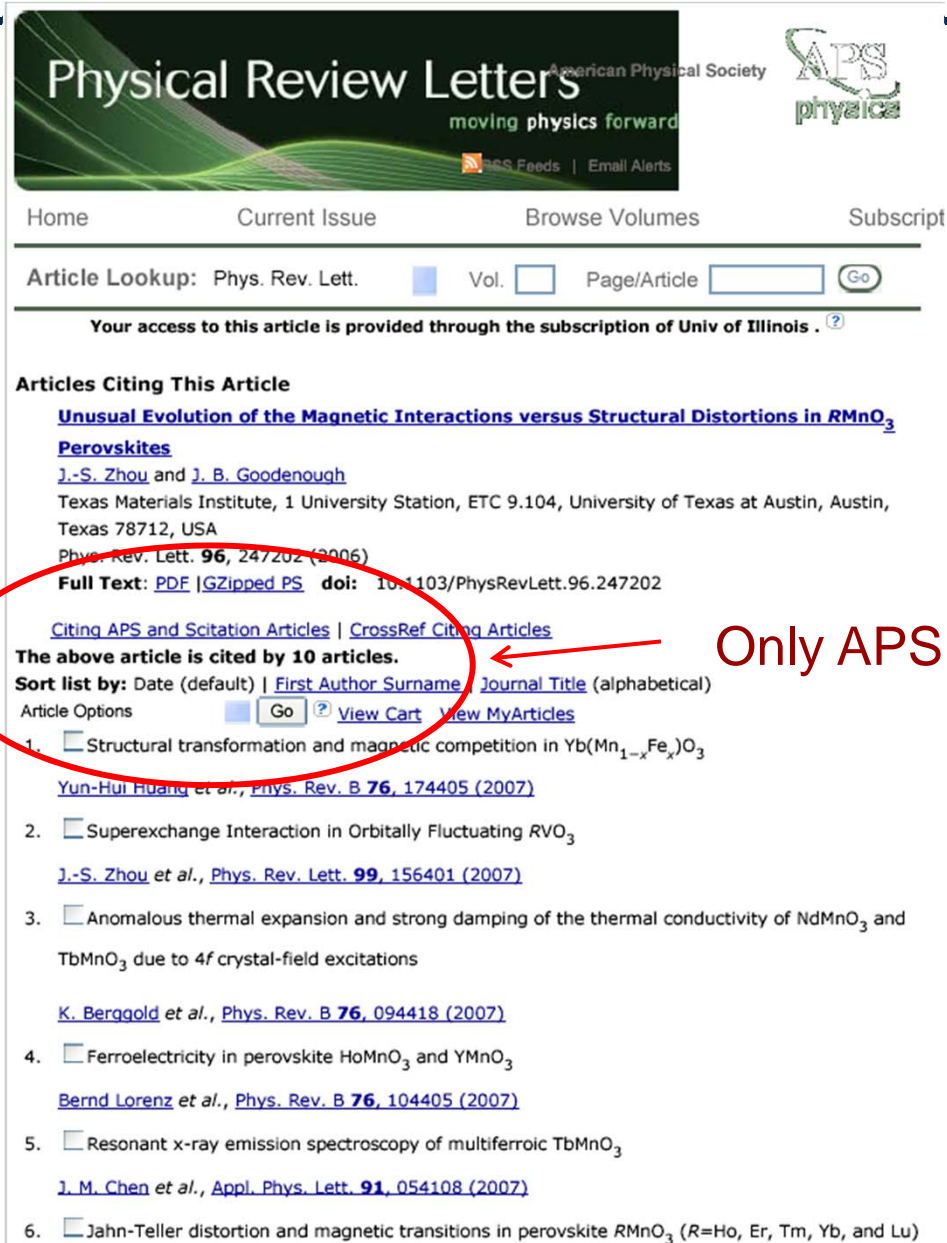
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1. J. B. Goodenough, [Phys. Rev.](#) **100**, 564 (1955).
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Title: Harmonic measure and winding of random conformal paths: A Coulomb gas perspective

Authors: [Bertrand Duplantier](#), [Ilia Binder](#)

Comments: 29 pages, 4 figures

Subjects: Statistical Mechanics (cond-mat.stat-mech)

We consider random conformally invariant paths in the complex plane (SLEs). Using the Coulomb gas method in conformal field theory, we rederive the mixed multifractal exponents associated with both the harmonic measure and winding (rotation or monodromy) near such critical curves, previously obtained by quantum gravity methods. The results also extend to the general cases of harmonic measure moments and winding of multiple paths in a star configuration.

[2] [arXiv:0802.2291](#) [[ps](#), [pdf](#), [other](#)]

Title: Competition between charge and spin order in the St-U-VS extended Hubbard model on the triangular lattice

Authors: [B. Davoudi](#), [S. R. Hassan](#), [A.-M. S. Tremblay](#)

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Subjects: Strongly Correlated Electrons (cond-mat.str-el)

Several new classes of compounds can be modeled in first approximation by electrons on the triangular lattice that interact through on-site repulsion U as well as nearest-neighbor repulsion V . This extended Hubbard model on a triangular lattice has been studied mostly in the strong coupling limit for only a few types of instabilities. Using the extended two-particle self consistent approach (ETPSC), that is valid at weak to intermediate coupling, we present an unbiased study of the density and interaction dependent crossover diagram for spin and charge density wave instabilities of the normal state at arbitrary wave vector. When U dominates over V and electron filling is large, instabilities are chiefly in the spin sector and are controlled mostly by

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Authors

1. Title: [3D framework structure of a new lithium thiophosphate](#)
Author(s): Kim Y, Arumugam N, Goodenough JB
Source: CHEMISTRY OF MATERIALS Volume: 20 Issue: 2 F
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2. Title: [Cathode materials: A personal perspective](#)
Author(s): Goodenough JB
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3. Title: [Correlation between the structure and the spin state in](#)
Author(s): Takami T, Zhou JS, Goodenough JB, et al.
Source: PHYSICAL REVIEW B Volume: 76 Issue: 14 Article N
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4. Title: [Structural transformation and magnetic competition in](#)
Author(s): Huang YH, Karppinen M, Imamura N, et al.
Source: PHYSICAL REVIEW B Volume: 76 Issue: 17 Article N
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Correlation between the structure and the spin state in $R_{1-x}Sr_xCoO_3$ (R=La, Pr, and Nd)

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Author(s): [Takami T](#) (Takami, T.), [Zhou JS](#) (Zhou, J. -S.), [Goodenough JB](#) (Goodenough, J. B.), [Ikuta H](#) (Ikuta, H.)

Source: PHYSICAL REVIEW B **Volume:** 76 **Issue:** 14 **Article Number:** 144116 **Published:** 2007

Times Cited: 0 **References:** 40

Abstract: The crystal structures of the perovskites $R_{1-x}Sr_xCoO_3$ (R=La, Pr, and Nd) with $0 \leq x \leq 0.3$ were refined from x-ray powder diffraction spectra with the Rietveld method. The result of the Rietveld analyses indicates the presence of three unequal Co-O bonds with different lengths for all samples, and we found that the difference between the two Co-O(2) bond lengths increases with temperature and Sr content. The intermediate-spin (IS) state of the Co^{3+} ion is Jahn-Teller (JT) active, and the structural data show that the concentration of Co^{3+} ions increases with temperature and Sr content. For the non-Sr-doped R=La sample, we observed only a small distortion at 90 K while the JT distortion was large at 300 K. With substitution of Sr for La, the distortion at 90 K was found to increase. This observation can be accounted for by the increased stability of the IS state with Sr doping. On the other hand, the distortion was small even at 300 K for nondoped $PrCoO_3$ and $NdCoO_3$, indicating that the low-spin state is stable up to higher temperatures than in $LaCoO_3$. The distortion at 90 K was larger for the Sr-doped samples as in the R=La system, but the Sr content at which the JT distortion started to increase significantly was larger than that in the R=La system.

Document Type: Article

Language: English

KeyWords Plus: MAGNETIC-PROPERTIES; ELECTRONIC-STRUCTURE; NEUTRON-DIFFRACTION; RCO_3R ; $RNiO_3R$; $LaCoO_3$; TRANSITION; PEROVSKITES; EU; SM

Addresses: Takami, T (reprint author), Univ Texas Austin, 1 Univ Stn, C2201, Austin, TX 78712 USA
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Example: O'Brian C* OR OBrian C*

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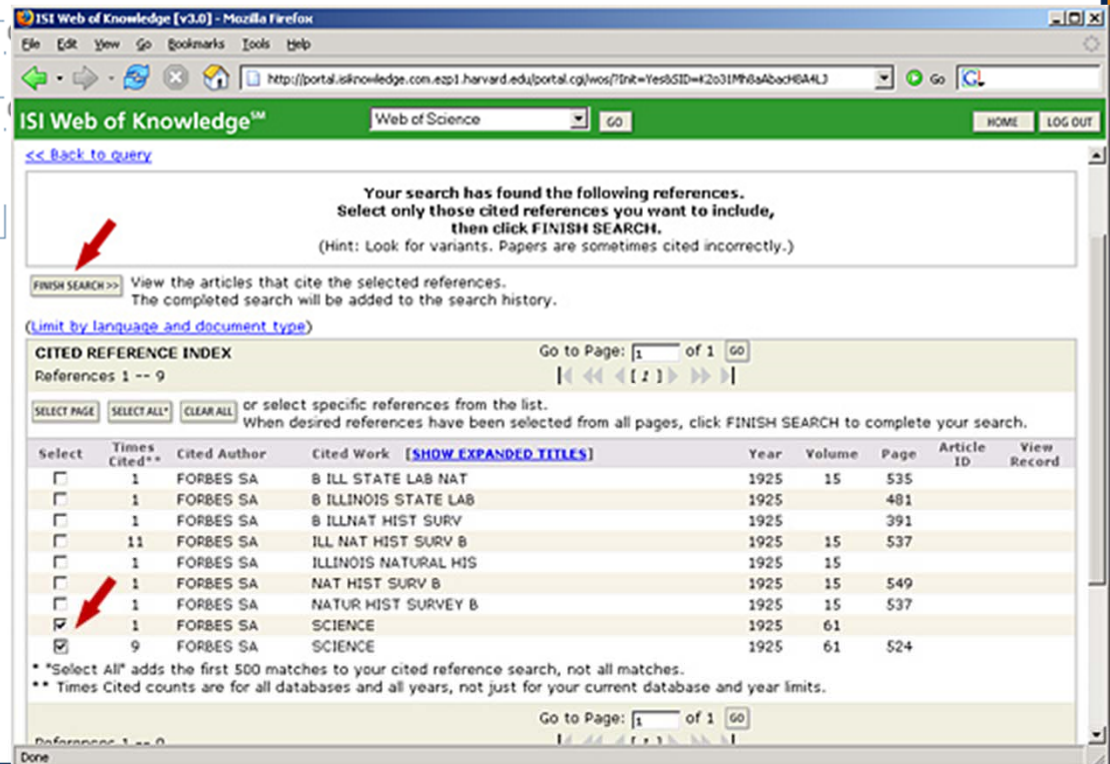
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Example: 1943 or 1943-1945

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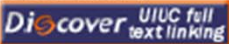



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Authors

Source Titles

Publication Years

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- Title: Ion exchange reactions of NaSbO₃ and morphotropic
Author(s): Nalbandyan VB, Avdeev M, Pospelov AA
Source: SOLID STATE SCIENCES Volume: 8 Issue: 12 Page Times Cited: 0
 Full Text
- Title: Disilver oxotellurate(VI), Ag₂TeO₄
Author(s): Klein W, Curda J, Peters EM, et al.
Source: ZEITSCHRIFT FUR ANORGANISCHE UND ALLGEME 2005
Times Cited: 8
 Full Text
- Title: On the crystal chemistry of oxoargentates and silverox
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Source: ZEITSCHRIFT FUR ANORGANISCHE UND ALLGEME
Published: 2004
Times Cited: 11
 Full Text
- Title: Synthesis and crystal structure of the novel Pb₅Sb₂Mn
Author(s): Abakumov AM, Rozova MG, Chizhov PS, et al.
Source: JOURNAL OF SOLID STATE CHEMISTRY Volume: 1:
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