Abstract Submitted for the MAR05 Meeting of The American Physical Society

Combining Dynamical Mean Field and Local Density Approximations in the study of GaMnAs J. MORENO, Univ. of North Dakota, M. A. MAJIDI, PAUL KENT, M. JARRELL, Univ. Cincinnati, R. S. FISHMAN, Oak Ridge Nat. Lab. — We use the band structure predicted by the Local Density Approximation as the starting point for a self consistent Dynamical Mean Field treatment of the ferromagnetic order in GaAs doped with Mn. Our model is an effective eight-band model including the heavy and light valence bands, the split-off valence band and the conduction band of the host material. The exchange interaction between the randomly distributed magnetic ions and the itinerant carriers is modeled using a modified double-exchange coupling. We compare our results with the ones predicted by more simplified models.

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Date submitted: 07 Dec 2004

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