Abstract Submitted for the MAR15 Meeting of The American Physical Society

Superconducting transition temperature in two-dimensional doped repulsive Hubbard model: DCA+ simulations with continuous momentum dependence¹ MI JIANG, Institute for Theoretical Physics, ETH Zurich, PETER STAAR, IBM Research - Zurich, THOMAS MAIER, Computer Science and Mathematics Division, Oak Ridge National Laboratory, THOMAS SCHULTHESS, Computer Science and Mathematics Division, Oak Ridge National Laboratory;Swiss National Supercomputing Center, ETH Zurich — DCA+ algorithm extends the dynamical cluster approximation (DCA) with continuous lattice self-energy to ensure better convergence with cluster size and delay the occurrence of the severe sign problem. This new algorithm enables a systematic investigation of the phase diagram of 2D Hubbard model relevant to the high temperature superconductors. We calculate the superconducting transition temperature T_c in the 2D repulsive Hubbard model on square lattice with nearest-neighbor hoppings for different doping levels, focussing on the intermediate correlation (U/t = 7) regime.

¹This research was carried out with resources of the Oak Ridge Leadership Computing Facility (OLCF), the Swiss National Supercomputing Center (CSCS), and the Center for Nanophase Materials Sciences (CNMS).

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Date submitted: 12 Nov 2014

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