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Supersolid phase of hardcore bosons on triangular lattice.

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We study properties of the supersolid phase observed for hardcore bosons on the triangular lattice near half-integer filling factor, and the phase diagram of the system at finite temperature. We find that the solid order is always of the $(2m, -m', -m')$ with $m$ changing discontinuously from positive to negative values at half-filling, in contrast with phases observed for Ising spins in transverse magnetic field. At finite temperature we find two intersecting second-order transition lines, one in the 3-state Potts universality class and the other of the Kosterlitz-Thouless type.

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Since the supersolid state of matter was introduced to physics nearly half a century ago and its theoretical feasibility was demonstrated,¹ there was a long history of experimental attempts to find it in Nature (mostly in $^4$He, see, e.g., Ref. 2) along with numerical simulations and theoretical predictions for models of interacting lattice bosons. Recent years have seen a renewed interest in this topic. On the one hand, lattice bosons are no longer restricted to $^4$He samples in the experimental point of view) models displaying a supersolid state persisting for smaller values of $t/V$. In Ref. [6] the system was thought to remain a disordered phase diagram is similar to that of Ref. [6], with the no- transition at $n=1/2$ and $t/V \approx 0.115$ and the stable supersolid state persisting for smaller values of $t/V$. In Ref. [6] the system was thought to remain a disordered superfluid for arbitrary $t/V$. The discrepancy can be attributed to known limitations of the GFMC method.

The model has been investigated in a series of recent papers, making use of advanced numerical techniques.⁸,⁹,¹⁰ The proposed zero-temperature phase diagram is similar to that of Ref. ⁶, with the notable addition of a quantum superfluid-supersolid phase transition at $n=1/2$ and $t/V \approx 0.115$ and the stable supersolid state persisting for smaller values of $t/V$. In Ref. ⁶ the system was thought to remain a disordered superfluid for arbitrary $t/V$. The discrepancy can be attributed to known limitations of the GFMC method.

Based on field-theoretic, exact diagonalization, and other arguments, Ref. ⁶ hints at the possibility of the
(m, 0, −m) density order in the ground state at n = 1/2 (state C). These considerations involved, in particular, an analogy between the properties of Eq. (2), and those of the Ising antiferromagnet on the triangular lattice, in the presence of a transverse magnetic field [11]. If true, there should exist quantum A − C and C − B phase transitions away from half-filling and three finite-temperature transitions of the Kosterlitz-Thouless (KT) type. Though Ref. [8] finds that the ground state is of the A or B type, it makes similar predictions for the finite temperature phase diagram at n = 1/2 which follow from the assumption that spontaneous symmetry breaking between A, B, and C, and their lattice translations is described by the six-clock model [13].

In what follows, we provide strong evidence that the supersolid state at half-filling is always of either the A or B type. Our data suggest that there is a discontinuous transition from A to B at μ = 3V similar to the I-order phase transition (driven by the large energy of the A − B domain walls). What makes it special is the exact particle-hole symmetry; structure factor, superfluid density, and energy remain continuous functions of μ through the transition line. For the supersolid A (or B) with the three-fold degenerate ground state, one expects to see the normal-superfluid KT and the solid-liquid 3-state Potts transitions, as temperature is increased. Moreover, the KT and Potts transitions are independent of each other and for n ≠ 1/2 intersect on the phase diagram. The failure of the mean-field description and analogies with the transverse-field Ising model to predict the supersolid structure at n = 1/2 can be traced back to the U(1)-symmetry of Eqs. (1) and (2), as noticed in Ref. [8]. For example, the (1, 0, −1) state can not be the true ground-state at finite t the limit of t/V → 0 simply because it does not respect the particle conservation law.

We use the worm-algorithm Monte Carlo scheme in the lattice path-integral representation [14] to simulate

\[
P(n^+) = \frac{\langle \sum_{k=1}^{N} \hat{n}_k e^{iQr_k} \rangle^2}{N^2} \quad (3)
\]

does not distinguish between supersolids A, B, C. We adopt the following strategy: for each system configuration, we compute the distribution of time-averaged occupation numbers, \( \bar{n}_k = T \beta^{-1} \int_0^\beta d\tau \bar{n}_k(\tau) \), and use it to determine the fraction of sites with \( \bar{n}_k > 1/2 \)

\[
n^+ = N^{-1} \sum_{k=1}^{N} \theta(\bar{n}_k - 1/2), \quad (4)
\]

where \( \theta(x) \) is the Heaviside function. A, B, C density structures correspond to \( n_A^+ = 2/3, n_B^+ = 0, \) and \( n_C^+ = 1/3 \). Finite systems are characterized by broad probability distributions \( P(n^+) \), and the formation of different solid orders can be seen as the development of sharp peaks, as the thermodynamic limit is approached.

In Fig. 2 we show the evolution of the \( P(n^+) \) distribution for the half-filled system at \( V/t = 10 \), i.e., close to the superfluid-supersolid transition point, estimated [8, 9, 10] at \( V/t \approx 8.5 \). The distribution is peaked at \( n^+ = 0 \) in the smallest system considered (\( L=6 \)), but, as the system size is increased, the weight is shifted toward the wings of the distribution. For \( L=18 \), there are already three peaks with comparable height. Finally, in the \( L=24 \) system we observe only two peaks corresponding to the supersolid phases A and B. Though the probability density between the peaks is still measurable, the dynamics of the algorithm becomes very slow; it typically takes millions of Monte Carlo sweeps, in order for the system to make a transition from the A to the B structure and vice versa. We have explicitly verified that configurations with \( n^+ \approx 2/3 \) and \( n^+ \approx 1/3 \) have density orders depicted as in Fig. 1I with a large contrast in density between sublattices. We have also checked that the

![FIG. 2: (Color online). Probability distributions \( P(n^+) \) for different system sizes and temperatures at \( \mu/V = 3 \) and \( t/V = 0.1 \).](image1)

![FIG. 3: (Color online). Probability distributions \( P(n^+) \) for different system sizes and temperatures at \( \mu/V = 3 \) and \( t/V = 0.05 \).](image2)
systems near half-filling, and a solid phase with algebraic correlations “sandwiched” between the solid and normal liquid phases. This prediction was made in Ref. 8 for $n = 1/2$. Since the ground state was found here to be only of the $A$ or $B$ type, and we do not see why domain wall energies between translated $A$ states are the same as between $A$ and $B$ states (in fact, the Landau theory prediction 8 is that $A$ and $B$ states phase separate and have different average densities even at $\mu = 3V$), the finite temperature phase diagram should instead feature the normal-superfluid KT and the liquid-solid 3-state Potts (for $n \neq 1/2$) transitions breaking $U(1)$ and translation symmetry respectively. At $n = 1/2$ we expect only one liquid-solid transition. An interesting question is whether transition lines simply intersect, or there are bicritical and tricritical points and I-order lines as observed for the similar model on the square lattice 13. We performed simulations for two representative cases, one for constant chemical potential $\mu/V = 2.74$ (or density $n \approx 0.44$), and the other for constant $t/V = 0.1$.

In Fig. 4 we show typical data for the KT transition between the solid and supersolid phases. The transition is smeared by logarithmic finite-size effects, but the critical temperature can be still determined with good accuracy by utilizing the well known renormalization flow and the universal jump of the superfluid density, $\rho_s$, at $T_c$. The data analysis is as follows: 10 we define $R = \pi \rho_s/2m^2T$ (where $m = 1/3t$ is the effective mass for the triangular lattice) and study the finite-size scaling of the data using KT renormalization group equations in the integral form

$$4 \ln(L_2/L_1) = \int_{R_2}^{R_1} \frac{dt}{t^2(\ln(t) - \kappa)} + t.$$  (5)

The microscopic (system size independent) parameter $\kappa$ is an analytic function of temperature, and the critical point corresponds to $R = 1$ at $\kappa = 1$. For $T < T_c$, the thermodynamic curve is defined by the equation

$$1/R + \ln R = \kappa(T),$$  (6)

with $\kappa = 1 + \kappa'(T_c - T)$. We use different pairs of system sizes in Eq. (5) to determine the $\kappa(T)$ curve, and obtain the location of the critical point from $\kappa(T_c) = 1$. The results are shown in the inset of Fig. 4. Data collapse and smooth analytic behavior of $\kappa(T)$ proves that the transition is indeed of the KT type. We used the same protocol and system sizes to determine other critical points.

In Fig. 5 we present our data for the transition into the state with the long-range density order. For the three-fold degenerate $B$ structure this transition is expected to be in the 3-state Potts universality class. The critical exponents are known exactly 17: $\nu = 5/6$, and $\beta = 1/9$. We thus perform the data collapse using $L^{2\beta}S_\rho = f(\delta L^{1/\nu})$ where $\delta = (T - T_c)/t$ and $T_c$ is the only fitting parameter. The result is shown in the inset of Fig. 5. This confirms the above-mentioned expectation, and establishes that there is only one transition to the solid phase (there are no visible finite-size effects below $T_c$).
Finally, we compute the phase diagram in the $(T/t, t/V)$ (at constant $\mu/V = 2.74$) and $(T/t, \mu/V)$ (at constant $t/V = 0.1$) planes and observe that KT and Potts transition lines form a simple cross for $n \neq 1/2$, i.e., the corresponding order parameter fields are not strongly interacting, see Fig. 6. The transition temperature to the superfluid and supersolid states in this part of the phase diagram is determined by the hopping amplitude. Within the statistical uncertainties of our calculation, KT and Potts transition temperatures cannot be distinguished at $\mu/V = 3$.

We did not see evidence for the algebraic solid state at $\mu = 3V$. The finite-size scaling for the supersolid-solid transition at $\mu = 3V$ is consistent with the 3-state Potts universality, though the data collapse is not as impressive as in Fig. 11 (the other alternative is the KT transition).

It is instructive to understand why the $(m, 0, -m)$ phase for the Hamiltonian (11) is not an obvious ground-state. At the mean-field level, $C$ has a better energy than $A$ or $B$. For the transverse-field Ising model (11) the $(1,0,-1)$ spin arrangement is obtained by orienting the middle spin along the magnetic field direction, i.e., putting it in the equal-amplitude superposition of up- and down-states. In bosonic language, it corresponds to the superposition of states with one or zero particles on a given site. Such a state can not be reconciled with the Hamiltonian (11) which conserves the particle number. Any non-integer average occupation number necessarily involves hopping transitions to the nearest neighbor sites. In the $(1,0,-1)$ structure the middle site is completely surrounded by the fully occupied or empty sites and thus can not be the ground state of the system even in the limit of $t/V \to \infty$. The problem appears to be inherently quantum with no obvious solution at the mean-field level.

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[7] Working with a finite population of random walkers in GFMC, has the effect of biasing estimates for the static structure factor, required to establish the presence of diagonal long-range order. Such a bias can in principle be eliminated by increasing the size of the population, but this is often impractical. See, for instance, M. Calandra Bonaura and S. Sorella, Phys. Rev. B 57, 11446 (1998).