Thermal fluctuations and quantum phase transition in antiferromagnetic Bose-Einstein condensates

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We present a method for investigating nonequilibrium dynamics of an ultracold system that is initially at thermal equilibrium. Our procedure is based on the classical fields approximation and appropriately prepared initial state for antiferromagnetic condensate with fixed magnetization. As an application of the method, we investigate the influence of thermal fluctuations on the quantum phase transition from an antiferromagnetic to phase separated ground state in a spin-1 Bose-Einstein condensate of ultracold atoms. We find that at temperatures significantly lower than the critical condensation temperature T_c the scaling law for the number of created spin defects remains intact.

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I. INTRODUCTION

The ground state phase diagram of an antiferromagnetic Bose-Einstein condensate was studied experimentally in the regime where spatial and spin degrees of freedom are decoupled [1], demonstrating a phase transition from an antiferromagnetic phase (where only the $m_F = \pm 1$ Zeeman components are populated) to a mixed broken axisymmetric phase (where all three Zeeman states can be populated). In this regime the system follows the predictions of the single spatial mode approximation. In other ranges of parameters, however, spatial separation of components can occur spontaneously, and spin domains may appear in the ground state of the system [2,3]. In this case the ground state is either the antiferromagnetic phase or is phase separated into two domains, and the two possibilities are divided by a critical point that is characterized by a critical magnetic field.

A system may become out of equilibrium when it is driven through the critical point due to the divergence of the relaxation time. If symmetry breaking occurs at the same time, this out-of-equilibrium process can produce various kinds of defects, depending on the dimensionality of the system and the form of the order parameter. The unified description of these phenomena is described by the Kibble-Zurek mechanism (KZM), which was studied in a number of physical systems, from the early Universe to ultracold atomic gases [4–11]. Among these, Bose-Einstein condensates of ultracold atoms offer realistic models of highly controllable and tunable systems [12,13]. Recent experiment with quasi-one-dimensional ultracold atoms confirmed the spontaneous creation of solitons via the Kibble-Zurek mechanism [14].

In recent papers [15,16], we demonstrated that the quantum phase transition from an antiferromagnetic to phase separated ground state in a spin-1 Bose-Einstein condensate of ultracold atoms exhibits scaling laws characteristic for systems displaying universal behavior. Phase separation leads to the formation of spin domains, with the number of domain walls depending on the quench time. Interestingly, the Kibble-Zurek scaling law was confirmed only for the dynamics close to the critical point. Further evolution led to postselection of domains, which gave rise to a second scaling law with a different exponent. The postselection was attributed to the conservation of an additional quantity, namely, the condensate magnetization.

In this paper, we develop a method for investigating nonequilibrium dynamics of an ultracold system that is initially

at thermal equilibrium. We apply this method to investigate the effect of nonzero temperature on the dynamics of the phase transition and the resulting scaling laws.

We consider this problem within the framework of the classical theory of a complex field with exact equations of motion being the Gross-Pitaevskii equations [17]. We explore the applicability of the method that was valid for single-component condensates. Nevertheless, extension of the method is not straightforward since there is an additional constant of motion-the magnetization. Conservation of the magnetization plays a crucial role in experiments and the state of the system depends on its value [1,18]. Indeed, classical fields or stochastic methods were applied for spinor condensates at nonzero temperature but for free magnetization [19]. Here we show how to prepare the initial state for antiferromagnetic spinor Bose-Einstein condensates that takes into account fixed magnetization. Unlike the spinor version of the stochastic Gross-Pitaevski equation [19], which is based on the existence of a thermalized reservoir, our method can be used to describe systems at temperatures much lower than the critical temperature of condensation.

The initial condition for the transition includes both the condensate and thermal atoms that introduce thermal fluctuations in the system. We sample the initial thermal equilibrium within the Bogoliubov approximation at a given temperature and for fixed total number of atoms and magnetization. Modes orthogonal to the condensate are thermally populated according to the Bogoliubov transformation. In our work, the field has to be interpreted not simply as the condensate wave function, but rather as the total matter field. We present both the results of single realizations of the field, which experimentally correspond to single experimental runs, and results of averaging over different initial states. We find that while the dynamics of the system can be altered by the thermal fluctuations, at relatively low temperatures the scaling law for the number of domains in the final state is intact.

II. THE MODEL

In the following, we consider a dilute, weakly interacting spin-1 Bose-Einstein condensate placed in a homogeneous magnetic field pointing along the z axis. For the sake of completeness, we recall the details of the model used previously in [15,16] to describe the dynamics of a condensate at zero

temperature within the truncated Wigner approximation. In the following section, we will generalize this approach to the case of a condensate at nonzero initial temperature.

We start with the Hamiltonian $H = H_0 + H_A$, where the symmetric (spin-independent) part is

$$H_0 = \sum_{j=-,0,+} \int dx \,\psi_j^{\dagger} \bigg(-\frac{\hbar^2}{2m} \nabla^2 + \frac{c_0}{2} \rho + V(x) \bigg) \psi_j. \quad (1)$$

Here the subscripts j = -,0,+ denote sublevels with magnetic quantum numbers along the magnetic field axis $m_f = -1,0,+1, m$ is the atomic mass, $\rho = \sum \rho_j = \sum \psi_j^{\dagger} \psi_j$ is the total atom density, and V(x) is the external potential. Here we restricted the model to one dimension, with the other degrees of freedom confined by a strong transverse potential with frequency ω_{\perp} . The spin-dependent part can be written as

$$H_{\rm A} = \int dx \left[\sum_{j} E_j \rho_j + \frac{c_2}{2} : \mathbf{F}^2 : \right], \tag{2}$$

where E_j are the Zeeman energy levels, the spin density is $\mathbf{F} = (\psi^{\dagger} F_x \psi, \psi^{\dagger} F_y \psi, \psi^{\dagger} F_z \psi)$, where $F_{x,y,z}$ are the spin-1 matrices and $\psi = (\psi_+, \psi_0, \psi_-)$. The spin-independent and spin-dependent interaction coefficients are given by $c_0 = 2\hbar\omega_{\perp}(2a_2 + a_0)/3 > 0$ and $c_2 = 2\hbar\omega_{\perp}(a_2 - a_0)/3 > 0$, where a_S is the *s*-wave scattering length for colliding atoms with total spin *S*. In the following analytic calculations we assume the incompressible regime where

$$c_0 \gg |c_2|,\tag{3}$$

which is a good approximation, e.g., in the case of 87 Rb or 23 Na spin-1 condensate.

The total number of atoms $N = \int \rho \, dx$ and magnetization $M = \int (\rho_+ - \rho_-) \, dx$ are conserved quantities. In reality, there are processes that can change both N and M, but they are relatively weak both in spin-1 ²³Na and ⁸⁷Rb condensates [18,20] and can be neglected on the time scales considered below.

The linear part of the Zeeman shifts E_j induces a homogeneous rotation of the spin vector around the direction of the magnetic field. Since the Hamiltonian is invariant with respect to such spin rotations, we consider only the effects of the quadratic Zeeman shift [2,3]. For sufficiently weak magnetic field we can approximate it by a positive energy shift of the $m_f = \pm 1$ sublevels $\delta = (E_+ + E_- - 2E_0)/2 \approx B^2 A$, where *B* is the magnetic field strength and $A = (g_I + g_J)^2 \mu_B^2/16E_{\text{HFS}}$, g_I and g_J are the gyromagnetic ratios of electron and nucleus, μ_B is the Bohr magneton, and E_{HFS} is the hyperfine energy splitting at zero magnetic field [2,3]. Finally, the spin-dependent Hamiltonian (2) becomes

$$H_{\rm A} = \int dx \left[AB^2(\rho_+ + \rho_-) + \frac{c_2}{2} : \mathbf{F}^2 : \right].$$
(4)

Except for the special cases $M = 0, \pm N$, the V(x) = 0 ground state phase diagram, shown in Fig. 1, contains three phases divided by two critical points at

$$B_1 = B_0 \frac{M}{\sqrt{2}N}, \qquad B_2 = B_0 \frac{1}{\sqrt{2}},$$
 (5)



FIG. 1. (Color online) Ground state phase diagram of an antiferromagnetic condensate for magnetization M = N/2. The green arrow indicates the direction of quench into a phase separated state considered in Sec. V.

where $B_0 = \sqrt{c_2 \rho / A}$ and ρ is the total density. The ground state can be (i) antiferromagnetic (2*C*) with $\psi = (\psi_+, 0, \psi_-)$ for $B < B_1$, (ii) phase separated into two domains of the 2*C* and $\psi = (0, \psi_0, 0)$ type (ρ_0) for $B \in (B_1, B_2)$, or (iii) phase separated into two domains of the ρ_0 and $\psi = (\psi_+, 0, 0)$ type (ρ_+) for $B > B_2$ [3]. What is more, the antiferromagnetic 2*C* state remains dynamically stable, i.e., it remains a local energy minimum up to a critical field $B_c > B_1$. Consequently, the system driven adiabatically from the 2*C* phase, across the phase boundary B_1 , and into the separated phase remains in the initial 2*C* state up to $B_c > B_1$ when the 2*C* state becomes dynamically unstable towards the phase separation.

III. MODELING THERMAL FLUCTUATIONS BY THE CLASSICAL FIELDS

A classical fields method is used for description of thermal effects of a single condensate [17]. Here we extend the applicability of the method to the system of spin-1 Bose-Einstein condensates with antiferromagnetic interactions.

In the method the condensed and noncondensed parts of the *j*th component are described by a complex function ψ_j . The description takes into account modes *k* with energies lower than the temperature:

$$\frac{\hbar^2 k_{\text{max}}^2}{2m} n_{\text{cut}} \simeq k_B T. \tag{6}$$

The relation between maximal momentum k_{max} and the temperature is not strictly defined and n_{cut} has no unique value [21]. In general, results of the classical field are cut-off dependent. Here, we assume that $n_{\text{cut}} = 1/3$.

In our simulations, we consider a lattice model for a classical field $\psi_j(x)$ with lattice spacing dx. We enclose the atomic field in the ring-shaped quasi-one-dimensional (1D) geometry with periodic boundary conditions at $\pm L/2$ and V(x) = 0. The total number of atoms is a constant of motion

$$dx \sum_{x,j} |\psi_j(x)|^2 = N,$$
(7)

as well as the total magnetization

$$dx \sum_{x} [|\psi_{+}(x)|^{2} - |\psi_{-}(x)|^{2}] = M.$$
(8)

The numerical evolution of the field is governed by the discretized counterpart of the Hamiltonian (1),(2).

The initial antiferromagnetic (2C) state at thermal equilibrium is prepared by employing Bogoliubov transformation of state 2C: $\psi_j = \tilde{\psi}_j + \delta \psi_j$ [16], with the constraint $\tilde{\psi}_0 =$ 0. Linearization of the Gross-Pitaevskii equation in small fluctuations $\delta \psi_j(t,x)$ around uniform background $\tilde{\psi}_j$ decouple fluctuations $\delta \psi_{\pm}$ from $\delta \psi_0$.

Fluctuations $\delta \psi_{\pm}$ are composed of phonon (*p*) and magnon (*m*) branches

$$\begin{pmatrix} \delta\psi_{+}\\ \delta\psi_{-} \end{pmatrix} = \begin{pmatrix} \sqrt{\rho_{+}}\\ \sqrt{\rho_{-}} \end{pmatrix} \sum_{k\neq 0}^{k_{\text{max}}} \left(b_{k}^{(p)} u_{k}^{(p)} e^{ikx} + b_{k}^{(p)*} v_{k}^{(p)*} e^{-ikx} \right) + \left(\frac{\sqrt{\rho_{-}}}{-\sqrt{\rho_{+}}} \right) \sum_{k\neq 0}^{k_{\text{max}}} \left(b_{k}^{(m)} u_{k}^{(m)} e^{ikx} + b_{k}^{(m)*} v_{k}^{(m)*} e^{-ikx} \right),$$

$$(9)$$

with quasiparticle energies

$$\epsilon_k^{(p)} = c_2 \rho \sqrt{\xi_s^2 k^2 [2(c_0/c_2) + \xi_s^2 k^2]},$$
(10)

$$\epsilon_k^{(m)} = c_2 \rho \sqrt{\xi_s^2 k^2 (8n_+ n_- + \xi_s^2 k^2)}.$$

Normalized modes satisfy

$$u_{k}^{(p)} \pm v_{k}^{(p)} = \left(\frac{\xi_{s}^{2}k^{2}}{2(c_{0}/c_{2}) + \xi_{s}^{2}k^{2}}\right)^{\pm 1/4},$$

$$u_{k}^{(m)} \pm v_{k}^{(m)} = \left(\frac{\xi_{s}^{2}k^{2}}{8n_{+}n_{-} + \xi_{s}^{2}k^{2}}\right)^{\pm 1/4},$$
(11)

where $n_{\pm} = \rho_{\pm}/\rho$. Here we use the spin healing length $\xi_s = \hbar/\sqrt{2mc_2\rho}$.

The small quadrupole mode fluctuations $\delta \psi_0$ [22] are given by

$$\delta\psi_0 = \sum_{k=0}^{k_{\text{max}}} \left(b_k^{(0)} u_k^{(0)} e^{ikx} + b_k^{(0)*} v_k^{(0)*} e^{-ikx} \right)$$
(12)

with their gapped spectrum for $b < b_c$,

$$\epsilon_k^{(0)} = c_2 \rho \sqrt{\left[\xi_s^2 k^2 + (1 - b^2)\right]^2 - \left(1 - b_c^2\right)^2}, \quad (13)$$

and the normalized eigenmodes

$$u_k^{(0)} \pm v_k^{(0)} = \left(\frac{(b_c^2 - b^2) + \xi_s^2 k^2}{2(1 - b_c^2) + (b_c^2 - b^2) + \xi_s^2 k^2}\right)^{\pm 1/4}.$$
 (14)

Here we use a rescaled dimensionless magnetic field

$$b = \frac{B}{B_0}.$$
 (15)

To generate the stochastic initial values of the classical field we proceed as follows. (i) For each realization, we generate the fluctuations $\delta \psi_j$ at temperature *T* by introducing thermal population of the Bogoliubov modes. In practice we generate complex numbers $b_k^{(x)}$ for $k \neq 0$ of $j = \pm$ components and for all *k* of j = 0 components according to the probability distribution

$$P(b_k^{(\mathbf{x})}) = \frac{1}{\pi} \frac{\epsilon_k^{(\mathbf{x})}}{k_B T} e^{-|b_k^{(\mathbf{x})}|^2/k_B T}.$$
 (16)

Here (x) denotes (m), (p), or (0). For a given realization, we build up fluctuations $\delta \psi_j$ according to Bogoliubov transformations (9) and (12). (ii) Then, we create the classical field with the constraint that the total atom number and magnetization

are fixed. The form of the field is $\psi_j = \frac{a_j}{\sqrt{L}} + \delta \psi_j$ with

$$a_{+} = \left(\frac{N - M - N_{\perp}}{2}\right)^{1/2},$$
 (17)

$$a_{-} = \left(\frac{N+M-N_{\perp}}{2}\right)^{1/2},$$
 (18)

where $N_{\perp} = \sum_{k,j} |\delta \psi_j|^2$ is the number of noncondensed atoms.

The ensemble of classical fields created in this way is the initial state for the time-dependent Gross-Pitaevskii (GP) equations

$$i\hbar\frac{\partial\psi_{0}}{\partial t} = \left(-\frac{\hbar^{2}\nabla^{2}}{2m} + c_{0}\rho\right)\psi_{0}$$

+ $c_{2}[(\rho_{+} + \rho_{-})\psi_{0} + 2\psi_{0}^{*}\psi_{+}\psi_{-}],$
 $i\hbar\frac{\partial\psi_{+}}{\partial t} = \left(-\frac{\hbar^{2}\nabla^{2}}{2m} + c_{0}\rho + AB^{2}\right)\psi_{+}$
+ $c_{2}[(\rho_{+} - \rho_{-})\psi_{+} + \rho_{0}\psi_{+} + \psi_{-}^{*}\psi_{0}^{2}],$
 $i\hbar\frac{\partial\psi_{-}}{\partial t} = \left(-\frac{\hbar^{2}\nabla^{2}}{2m} + c_{0}\rho + AB^{2}\right)\psi_{-}$
+ $c_{2}[(\rho_{-} - \rho_{+})\psi_{-} + \rho_{0}\psi_{-} + \psi_{+}^{*}\psi_{0}^{2}].$ (19)

One of the advantages of the method is the ability to calculate various correlation functions in a straightforward way, taking averages over many realizations of the stochastic fields.

Application of the method is justified for a very low temperature only since Bogoliubov transformation is the approximate solution. Moreover, very early evolution of the GP equations can display some transient effects due to the fact that the Bogoliubov transformation used in the sampling does not produce an exactly stationary distribution. We have checked if such transients occur for parameters used in our simulations and found them to be marginal for the quantities we are interested in.

IV. VALIDITY OF THE BOGOLIUBOV TRANSFORMATION

The dynamics with Gross-Pitaevskii equations leads the two-component antiferromagnetic initial state to equilibrium state for long enough times. The equilibration given by this coupled set of nonlinear equations is surprising because we know they may lead to the coherent off-equilibrium spinmixing dynamics. Indeed, it was shown by using mean-field theory and adapting the single-spatial-mode approximation, that the condensate dynamics is well described by a nonrigid pendulum and displays a variety of periodic oscillations [18,23]. Fortunately, the period of oscillations depends on the initial fraction of atoms in the $m_F = 0$ component and is almost infinite in our case. Therefore, the time scale associated with spin-multimode dynamics is much shorter and allows for equilibration of the system. The relaxation observed in the system is quite intriguing, in particular, in the context of prethermalization phenomena [24].



FIG. 2. (Color online) Momentum distribution of noncondensed fields $|\delta \psi_j(k)|^2$ for $m_F = 1$ (left panel), $m_F = 0$ (center panel), and $m_F = -1$ (right panel). Green: initial distribution; red: after 100 s of equilibration with GP equations; black: Bogoliubov transformation Eqs. (9) and (12). Here $N = 10^7$, M = N/2, b = 0, with length $L = 200 \ \mu m$ and $\omega_{\perp} = 2\pi \times 1000 \ \text{Hz}$; other parameters as for sodium.

In the low temperature limit, the Bogoliubov theory well describes the equilibrium state of the system. The comparison of the result of equilibration with GP equations to the Bogoliubov theory can be treated as an independent test of the last one. To this end, we checked the validity of the Bogoliubov transformation, (9) and (12), as well as the quasiparticle energies, (10) and (13). We start numerical calculations with randomly chosen initial fluctuations with flat distribution in momentum space, see green points in Fig. 2, and superimposed constraints of given norm N and magnetization M. Then we let the initial state evolve with Gross-Pitaevskii equations for a transient time (100 s for results presented in Figs. 2 and 3). Next, we compare the achieved distribution of the noncondensed field in momentum space $\delta \psi_i (k \neq 0)$ to the predictions of the Bogoliubov theory. The comparison shown in Fig. 2 is satisfactory. Analysis in the frequency domain allows for a test of the validity of the elementary excitation picture (see the density plots in Fig. 3). One can easily recognize the magnon and phonon branches for $m_F = \pm 1$ and the quadrupole branch for $m_F = 0$. Once again, the comparison with the Bogoliubov theory (marked by black dashed lines) is satisfactory and shows that the quasiparticle energies (10) and (13) are recovered by equilibration applied to the initial state.

V. THE KIBBLE-ZUREK MECHANISM

In [15,16] we investigated the phase transition from the antiferromagnetic to the phase separated state in a spin-1 Bose-

Einstein condensate. This continuous phase transition is driven by the change of the magnitude of the applied magnetic field. Due to the spatial symmetry breaking in the phase separated state, the transition is accompanied by the creation of multiple defects in the form of spin-domain walls, with the number of domains dependent on the quench time. The concept of the KZM relies on the fact that the system does not follow the ground state exactly in the vicinity of the critical point due to the divergence of the relaxation time. The dynamics of the system cease to be adiabatic at $t \simeq -\hat{t}$ (here we choose t = 0in the first critical point), when the relaxation time becomes comparable to the inverse quench rate

$$\hat{\tau}_{\rm rel} \approx |\hat{\varepsilon}/\hat{\varepsilon}|,$$
 (20)

where $\varepsilon(t) = B - B_c \sim t/\tau_Q$ is the distance of the system from the critical point. At this moment, the fluctuations approximately freeze, until the relaxation time becomes short enough again. After crossing the critical point, distant parts of the system choose to break the symmetry in different ways, which leads to the appearance of multiple defects in the form of domain walls between domains of 2C and ρ_0 phases. The average number of domains is related to the correlation length $\hat{\xi}$ at the freeze-out time $\hat{t} \sim \tau_0^{z\nu/(1+z\nu)}$ [6,25],

$$N_{\rm d} = L/\hat{\xi} \sim \tau_{\rm O}^{-\nu/(1+z\nu)},$$
 (21)

where z and v are the critical exponents determined by the scaling of the relaxation time $\tau_{rel} \sim |\varepsilon|^{-zv}$ and excitation spectrum $\omega \sim |k|^{z}$, with z = 1 in the superfluid.



FIG. 3. (Color online) Spectral density $|\psi_j(k,\omega)|^2$, for $m_F = 1$ (left panel), $m_F = 0$ (center panel), $m_F = -1$ (right panel) after 100 s of evolution with GP equations. Black dashed lines: quasiparticle energies (10) and (13) given by the Bogoliubov theory. Parameters of the simulation are the same as in Fig. 2.

Interestingly, the Kibble-Zurek scaling law gives correct predictions only for the dynamics close to the critical point. Further on, the postselection of domains was observed, which gave rise to a second scaling law with a different exponent. The postselection was attributed to the conservation of an additional quantity, namely, the condensate magnetization.

The analytical and numerical calculations of [15,16] were carried out within the zero-temperature limit of the truncated Wigner approximation. In this section we use the method of Sec. III to estimate the influence of nonzero temperature on the dynamics of the phase transition and the resulting scaling laws. We find that while the dynamics of the system can be altered by the thermal fluctuations, at relatively low temperatures the scaling law for the number of domains in the final state is intact.

We now describe in detail the scenario of the experiment. The antiferromagnetic spin-1 condensate is trapped in a ringshaped quasi-1D trap with strong transverse confinement and the circumference length L. The magnetic field is initially switched off, and the atoms are prepared in the homogeneous antiferromagnetic (2C) ground state with magnetization set to M = N/2. To investigate KZM we increase B linearly as

$$B(t) = B_0 \frac{t}{\tau_Q},\tag{22}$$

to drive the system through the two phase transitions into a phase separated state. Then, at $t > \tau_Q$ the magnetic field is kept constant at the level $B = B_0$. As described in Sec. II, the ground state of the system becomes separated into the 2C and the ρ_0 phase at $B = B_1$, but the initial 2C state remains metastable until the point $B = B_c$. At this point, the system is expected to undergo phase transition accompanied by spatial symmetry breaking. As described above, due to the finite quench time the phase transition has a nonequilibrium character, and multiple spin domains can develop in the system, instead of two as the form of the ground state would suggest. Further, at the second critical point $B = B_2$, there is no symmetry breaking accompanying the phase transition and the spin-domain landscape remains intact. The mean-field critical exponents of the symmetry-breaking phase transition are z = 1 and v = 1/2, which according to the formula (21) gives the scaling law for the number of domains as $N_d \sim \tau_{\rm O}^{1/3}$. However, as shown in [15,16], this prediction is correct only for the number of domain seeds formed close to the critical point. When the domains become fully formed, their number is decreased in the postselection process, which is due to the existence of an upper limit for the number of domains in a system with conserved magnetization M.

Here, we investigate the influence of the finite temperature on the process described above. In Fig. 4 we show the evolution of the density of atoms in the initially unoccupied $m_f = 0$ state during the process of domain formation. The top figure shows a single realization of the zero-temperature truncated Wigner simulation, which can be interpreted as a result of a single experiment. For comparison, an analogous result in the case of finite temperature, obtained using the method described in Sec. III, is shown in the figure below. While this particular example corresponds to a relatively short quench time, when the postselection process is not very effective, it is visible that some of the initial fluctuations merge to form a single domain



FIG. 4. (Color online) Formation of spin domains through a modified Kibble-Zurek mechanism in a ring-shaped 1D geometry with ring length $L = 200 \ \mu \text{m}$ and $\omega_{\perp} = 2\pi \times 1000 \text{ Hz}$, for $N = 10^6$ atoms. The evolution of the density of the $m_f = 0$ component $|\psi_0|^2$ in a single realization of the experiment is shown. The top and bottom panels correspond to the zero-temperature and finite-temperature results. The quench time is taken as $\tau_{\rm Q} = 100$ ms. Total condensed fraction at t = 0 is $N_{k=0}/N = 0.89$.

instead of two. The effect of the finite temperature can be seen as the picture of the dynamics being more "fuzzy" in the figure below, but the number of created domains and their properties seem to be unaffected due to the relatively low temperature.

This conclusion is further confirmed in Fig. 5, where we show the results of systematic averaging of the number of created domains over many realizations of the initial distribution. The two sets of points showing the result for T = 0 and finite temperature overlap except for the very long times, where some decrease of the number of domains in the T > 0 case can be observed. Close inspection of the dynamics leads us to account this slight decrease (on average by less than one domain) to the lower number of domains created at the symmetry phase transition when the initial state already contains thermal fluctuations. In particular, the finite temperature does not significantly affect the $N_d \sim \tau_Q^{2/3}$ scaling law predicted to result from the domain postselection process. This turns out to be the case for any temperature investigated



FIG. 5. (Color online) Comparison of the averaged number of spin domains N_d after the quench as a function of the quench time for $N = 10^6$ atoms. The points are results of numerical simulations averaged over 100 runs. The square and circle points correspond to zero-temperature and finite-temperature results. The scale is logarithmic on both axes and N_d is decreased by 2 to account for the ground state phase separation into two domains. The dashed line is the fit to the power law with scaling exponent $n_d = -0.73 \pm 0.04$. The grid size is 2^8 and the cut-off is equal to 1/3.

by our method based on the Bogoliubov transformation. We note that the scaling law may be affected by thermal effects at temperatures higher than the ones achievable using the current method. This will be the topic of a future study.

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VI. CONCLUSIONS

In summary, we developed a method for investigating nonequilibrium dynamics of an ultracold system that is initially at thermal equilibrium. Our procedure is based on the classical fields approximation with an appropriately prepared initial state. We described in detail how to model thermal fluctuations using the Bogoliubov transformation, and demonstrated its validity by performing dynamical equilibration of an antiferromagnetic initial state with given total atom number and magnetization. We studied the effect of the nonzero temperature on the scaling law for the number of domains created via the Kibble-Zurek mechanism in antiferromagnetic spinor condensates. The effect of the finite temperature is visible on the evolution of the gas density. We find the density to be more "fuzzy" but the number of created domains and their properties rather unaffected by the temperature. This result in a relatively low temperature shows the strength of the conservation law, namely, the conservation of magnetization in our system that enforces the scaling law. In addition to the main result, we observed relaxation of the antiferromagnetic initial state to a thermallike equilibrium. This intriguing feature provides an interesting direction for future work.

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