Topological thermalization via vortex formation in ultrafast quenches

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We investigate the thermalization of a two-component scalar field across a second-order phase transition under extremely fast quenches. We find that vortices start developing once the thermal bath sets the control parameter to its final value in the nonsymmetric phase. Specifically, we find that vortices emerge as the fluctuating field relaxes and departs macroscopically from its symmetric configuration. The density of primordial vortices at the relaxation time is a decreasing function of the final temperature of the quench. Subsequently, vortices and antivortices annihilate at a rate that eventually determines the total thermalization time. This rate decreases if the theory contains a discrete anisotropy term, which otherwise leaves the primordial vortex density unaffected. Our results thus establish a link between the topological processes involved in the vortex dynamics and the delay in the thermalization of the system.

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I. INTRODUCTION

The network of topological defects left behind a phase transition is a ubiquitous phenomenon in nature which encodes valuable information about the underlying dynamics and the corresponding field theory in a number of systems [1-23]. Precisely in this context, Kibble [24,25] and Zurek [26] performed a seminal attempt to relate the critical dynamics with the spatial density of topological defects, which is commonly known as the Kibble-Zurek (KZ) mechanism. Their original arguments go as follows. In a finite-temperature quench, as the transition proceeds, the actual correlation length of the order-parameter distribution cannot forever follow its nominal equilibrium value, which diverges at the critical temperature. This dropout can be related to causality, since the speed of signal propagation remains finite. As a result, the correlation length gets frozen at some value which determines the size of the domains on which the order parameter takes approximately uniform values on the ground-state manifold (i.e., vacuum manifold). The random choice of different values at different locations of the sample is at the origin of spontaneous symmetry breaking. According to Kibble's argument [24], the evolution of the order parameter at the interface between two domains is governed by the so-called geodesic rule. That is, the order parameter interpolates through each interface following the shortest path on the vacuum manifold. In turn, this determines whether topological defects form. From this reasoning, it follows that the value of the correlation length at the freezing time determines the typical distance between topological defects. The KZ rationale has been applied to interpret the formation of topological defects in several condensed-matter setups, which has revealed intriguing departures from the postulated KZ scaling in the case of fast quenches [20-23].

Here, we investigate numerically the formation of vortices in a two-component classical scalar field under extremelyfast-quenching conditions, which represents a limit case of the usual KZ scenario. This implies that the vortices emerge in a process that is entirely driven by the dynamics of the system after the quench. In particular, the order-parameter correlations evolve according to its diffusive fluctuating dynamics at the final temperature of the quench, rather than being imprinted by its critical behavior just before the freezing. Diffusion is caused by thermal fluctuations which, in contrast with previous treatments, we simulate in a self-consistent manner throughout the quenches. Ultimately, we find that the process of formation and annihilation of vortices results in an effective "topological delay" in the thermalization of the system. In addition, we investigate the impact of a discrete sixfold anisotropy in the system, as this applies to some condensed-matter setups of interest.

The paper is organized as follows. In Sec. II we describe the fundamentals of the model and explain our approach. In Sec. III we present the results of our study. We summarize our conclusions in Sec. IV.

II. PRELIMINARIES

A. The model

We perform numerical simulations upon an effective Lagrangian model of a scalar field order parameter, Q, in which temperature fluctuations are incorporated through a Langevin term in the equation of motion. Specifically, we consider a two-component scalar order parameter, $Q = (Q_1, Q_2) =$ $(Q \cos \phi, Q \sin \phi)$, with equations of motion

$$\frac{\partial \mathcal{L}}{\partial Q_i} - \nabla \frac{\partial \mathcal{L}}{\partial \nabla Q_i} = -\frac{\partial \mathcal{R}}{\partial \dot{Q}_i} + f_i.$$
 (1)

Here \mathcal{L} is the effective Lagrangian, \mathcal{R} is the dissipative function, and f_i are the components of a stochastic Langevin force such that $\langle f_i(\mathbf{r}, t) f_j(\mathbf{r}', t') \rangle = 2\gamma T \delta_{ij} \delta(\mathbf{r} - \mathbf{r}') \delta(t - t')$, with T being the temperature and γ the damping coefficient [27]. In this way, we effectively consider Gaussian fluctuations of \mathcal{Q} satisfying the fluctuation-dissipation theorem at temperature T [28]. For the sake of consistency, the effective Lagrangian is evaluated at the same temperature.

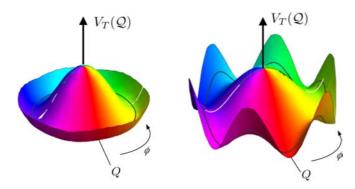


FIG. 1. Sketch of the total effective potential of the order parameter in configuration space, $V_T(Q) = V(Q) + V_6(Q)$, below the critical temperature. The field values at the minimum of $V_T(Q)$ form the vacuum manifold. Left: potential in the weak-anisotropy regime. Right: potential in the strong-anisotropy regime.

This is in contrast with previous approaches (cf. Refs. [29,30]) in which the amplitude of the noise term is kept constant and irrespective of the value of the temperature which enters the effective Lagrangian. Further, the noise amplitude is chosen there small enough so as to hardly affect the dynamics of the phase transition. We will see later that our requirement of consistency in the Langevin term is crucial in the evolution of the transition.

In the following, we consider an effective Lagrangian of the form

$$\mathcal{L} = V_T(\mathcal{Q}) + \frac{g}{2} [(\nabla \mathcal{Q})^2 + \mathcal{Q}^2 (\nabla \phi)^2], \qquad (2)$$

where the total potential V_T contains a U(1)-symmetric term, V(Q), and a sixfold anisotropy term, $V_6(Q)$, $V_T(Q) = V(Q) + V_6(Q)$,

$$V(Q) = \frac{a}{2}Q^2 + \frac{b}{4}Q^4 + \frac{c}{6}Q^6,$$
(3)

$$V_6(\mathcal{Q}) = \frac{c'}{6} \left[\left(Q_1^2 - 3Q_2^2 \right) Q_1 \right]^2 = \frac{c'}{6} Q^6 \cos^2(3\phi).$$
(4)

As is customary, we define $a = a_0 \epsilon$ as the control parameter with $\epsilon = (T - T_c)/T_c$ being the reduced temperature; its temperature dependence results from the effective coupling of

the system to the thermal bath. The coefficients b, c, g > 0 are all constant. The total effective potential according to these definitions is sketched in Fig. 1. The critical behavior of this model belongs to the XY universality class. However, the sixfold anisotropy term, $V_6(Q)$, can be tuned to describe either U(1)-symmetric systems such as superfluids or \mathbb{Z}_6 -symmetric ones such as hexagonal multiferroic manganites [31–33]. In the latter case, although the symmetry that is initially broken below T_c is that of the U(1) group, the subsequent evolution as well as the final structure of the vortices is generally affected by the \mathbb{Z}_6 anisotropy term for sufficiently large values of c'.

In static equilibrium, the uniform order parameter for which the free energy presents a minimum value satisfies

$$\{a + bQ^2 + [c + c'\cos^2(3\phi)]Q^4\}Q = 0,$$
 (5)

$$Q^6 \sin(6\phi) = 0. \tag{6}$$

The only solution above T_c (a > 0) is $Q_1 = Q_2 = 0$. Below T_c (a < 0), the above equation presents 12 possible solutions with $Q = Q_0 \simeq \sqrt{|a|/b}$, $\phi_n = n\pi/12$ (n = 0, ..., 11). However, the actual minima of the energy correspond to either $\phi_n = n\pi/3$ if c' < 0, or $\phi_n = (2n + 1)\pi/6$ if c' > 0, where n = 0, ..., 5, unless higher-order terms are included [32]. Thus, the minimum energy state of the order parameter is sixfold degenerate. Hence, the choice of any of those values by the order parameter breaks spontaneously the symmetry of the system whose phase is then said to be nonsymmetric.

Regarding the dynamics, we neglect inertial terms and restrict ourselves to the overdamped dynamics of Q introducing the dissipative function \mathcal{R} [27],

$$\mathcal{R} = \frac{\gamma}{2} (\dot{Q}_1^2 + \dot{Q}_2^2) = \frac{\gamma}{2} (\dot{Q}^2 + Q^2 \dot{\phi}^2), \tag{7}$$

in accord with the Langevin force [28].

All in all, the time evolution of the system is described by the stochastic diffusion equation (1) which contains a radial force f_Q acting upon the amplitude of the order parameter, and a tangential force f_{ϕ} upon its phase. For future purposes, it is convenient to distinguish three contributions to these forces as follows:

$$-\gamma \partial_t Q \equiv f_Q = \underline{aQ + bQ^3 + [c + c'\cos^2(3\phi)]Q^5} \underbrace{-g[\nabla^2 Q - Q(\nabla\phi)^2]}_{\text{Therefore}} \underbrace{-f_1\cos\phi - f_2\sin\phi}_{\text{Therefore}},\tag{8}$$

$$-\gamma \partial_t \phi \equiv f_{\phi} = \underbrace{-c' \sin(6\phi) Q^4/6}_{-g Q^{-2} \nabla(Q^2 \nabla \phi)} \underbrace{+Q^{-1}(f_1 \sin \phi - f_2 \cos \phi)}_{-g Q^{-2} \nabla(Q^2 \nabla \phi)}.$$
(9)

tangential restoring force tangential tension force tangential stochastic force

Thus, we identify a *restoring force* associated to the effective potential, a *tension force* associated to gradient terms, and a *stochastic force* originated from the Langevin term.

since equations can be obtained for the amplitude and phase zero modes, i.e., for uniform values of Q and ϕ , respectively.

In the symmetric phase, above T_c , the perturbations of the order parameter around its stable point (0,0) can be written such that $(Q_1, Q_2) = (0, 0) + (q_1, q_2)$. Thus, the effective Lagrangian associated to the Gaussian fluctuations is

B. Correlation lengths and relaxation times in the mean-field approximation

In this section we outline the definition of the correlation lengths and relaxation times in the mean-field approximation,

$$\delta \mathcal{L} = \frac{g}{2} \left[\bar{\xi}^{-2} \left(q_1^2 + q_2^2 \right) + (\nabla q_1)^2 + (\nabla q_2)^2 \right], \quad (10)$$

where $\bar{\xi} = (g/a)^{1/2}$ is the correlation length, common to the q_1 and q_2 components. Likewise, linearization of the equations of motion in the symmetric phase around Q = 0, neglecting the Langevin force and the tensions, yields

$$-\partial_t q_{1,2} = (|a|/\gamma)q_{1,2}, \tag{11}$$

from which we identify $\overline{\tau} = \gamma/|a|$ with a relaxation time common to q_1 and q_2 .

Below T_c , once in the nonsymmetric phase, it is more convenient to write Q in terms of Q and ϕ instead. Hence, the value of the order parameter around any of the six minima of the effective potential can be written now as $(Q, \phi) =$ $(Q_0, \phi_n) + (q, \varphi)$. In this case, the effective Lagrangian for the Gaussian fluctuations reads

$$\delta \mathcal{L} = \frac{g}{2} \{ \left[4\bar{\xi}^{-2}q^2 + (\nabla q)^2 \right] + Q_0^2 \left[\bar{\xi}_6^{-2}\varphi^2 + (\nabla \varphi)^2 \right] \}.$$
(12)

It is plain that, in addition to the correlation length of the amplitude zero mode, $\bar{\xi}$, there appears a second correlation length $\bar{\xi}_6 = [g/(3|c'|Q_0^4))^{1/2}$ associated to the fluctuations of the phase zero mode, φ . Again, linearization around (Q_0, ϕ_n) as for Eq. (12) yields the following equations for the fluctuations of the amplitude and the phase modes, respectively:

$$-\partial_t q = \bar{\tau}^{-1} q, \tag{13}$$

$$-\partial_t \varphi = (3|c'|Q_0^4/\gamma)\varphi, \qquad (14)$$

where both the Langevin force and the tension forces have been neglected. Again, φ presents a second relaxation time, $\bar{\tau}_6 = \gamma/(3|c'|Q_0^4)$.

C. Numerical simulations

We perform numerical simulations on a cubic sample, considering two different situations. In the first place, we consider that the system is effectively U(1)-symmetric, and refer to this situation as weak-anisotropy regime. That is, we choose the parameters in the potential V_T such that the anisotropic term becomes irrelevant, i.e., $V_T(\mathcal{Q}) \approx V(Q)$. In particular, for the simulations in Sec. III A, we take for the side lengths, $L_x = L_y \approx 400l$, $L_z \approx 120l$, where l is the lattice spacing. The dynamics of the order parameter Q is governed by a discretized version of Eqs. (8) and (9). The numerical values of the parameters in those equations are $a_0 = 1, b = 2$, $c = 0, c' = 2/3, g = 1, T_c = 0.0025$, which are chosen such that $a_0^{-1/2}$ becomes the unit of length scale and, at temperature $T_c/2, \epsilon = -0.5, \bar{\xi}(T_c/2) = 1/\sqrt{2}, \text{ and } \bar{\tau}(T_c/2) = 0.2 \text{ for } \gamma =$ $0.1a_0^{1/2}$. In terms of the correlation lengths and relaxation times of the mean-field approximation, this implies also that $\bar{\xi} \ll \bar{\xi}_6$, $\bar{\tau} \ll \bar{\tau}_6$ in the weak-anisotropy regime (cf. Table I below).

As for the situation in which $V_6(Q)$ becomes relevant, that we refer to as *strong-anisotropy regime*, the coefficient c' is enlarged so that the phase zero mode gets massive and the relationship between the mean-field correlation lengths and relaxation times turns into $\bar{\xi} > \bar{\xi}_6$, $\bar{\tau} > \bar{\tau}_6$. In particular, for the simulations in Sec. III B, we take for the side lengths $L_x = L_y = 1314l$, $L_z = 43l$. The numerical values of the rest of the parameters are as in the weak-anisotropy regime, except for the value of c' that is taken as c' = 128/3. Simulations in

TABLE I. Compilation of the results of the numerical simulations for different final temperatures, ϵ_f , in the weak-anisotropy regime. Lengths and times are given in units of $a_0^{-1/2}$, whereas the values for n_0 are in units of a_0 .

ϵ_{f}	-0.3	-0.4	-0.5	-0.6	-0.7	-0.8
$n_0[a_0] \ \xi_0[a_0^{-1/2}]$	0.0021	0.0027	0.0036	0.0041	0.0042	0.0054
	21.77	19.21	16.61	15.56	15.35	13.64
ર્ક	0.91	0.79	0.71	0.65	0.60	0.56
ર્ક્6	4.71	3.54	2.83	2.36	2.02	1.77
$ \begin{aligned} \overline{\tau_d[a_0^{-1/2}]} \\ \tau_0 \\ \tau_1 \end{aligned} $	1.4	1.1	0.88	0.77	0.68	0.62
	1.9	1.4	1.1	0.97	0.82	0.71
	2.61	1.98	1.59	1.37	1.19	1.05
$ar{ au} \ ar{ au}_6$	0.33	0.25	0.20	0.17	0.14	0.13
	2.2	1.3	0.80	0.56	0.41	0.31

the strong-anisotropy regime are performed for a single final temperature, $\epsilon_f = -0.3$, for which $\bar{\xi} = 0.71$, $\bar{\xi}_6 = 0.59$.

Since the vortex radius scales approximately with the minimum correlation length of the mean-field approximation, the lattice spacing *l* is adapted in either case to each temperature according to the formula $l = \min(\sqrt{2\xi}, \xi_6)/4$. It is worth mentioning that, with the numerical values chosen for the parameters, the Ginzburg region reduces to $\epsilon \in [-3.96 \times 10^{-10}, 3.96 \times 10^{-10}]$, which is negligible in all our simulations. Therefore, effects related to the strongly nonperturbative dynamics of the order parameter within this region can be fairly discarded. Lastly, the boundary conditions imposed upon Q are of the kind of no-flux boundary conditions, $\mathbf{n} \cdot \nabla Q_{1,2}|_{\partial\Omega} = 0$, which physically means that the polarization vector remains fixed at the boundary $\partial\Omega$, $\mathbf{n} \perp \partial\Omega$.

In all the simulations we consider ultrafast quenches from the initial temperature $T_i > T_c$ in the symmetric phase, to the final temperature $T_f < T_c$, assuming a uniform rate $(T_f - T_c)$ $T_i)/\tau_{\rm cool}$ with $\tau_{\rm cool}$ being the cooling time. Ultrafast quenches are defined by the inequality $\tau_{cool}(T_c - T_f)/(T_i - T_f) \ll \tau_0$, where the term on the left-hand side of this inequality is the time lapse corresponding to the temperature interval $[T_c, T_f]$, and τ_0 is the vortex formation time measured from the time the quenching passes through T_c . This means that the time spent by the system during the quench below T_c is negligible in comparison to the vortex formation time τ_0 . Also, it implies that its dynamics is independent of the initial temperature. It is important to note that in our simulations the stochastic forces act on the system all the way through from the start. Finally, since the dynamics of Q from Eqs. (8) and (9) is determined by the ratio T/γ , we fix the value of the diffusion coefficient at $\gamma = 0.1 a_0^{1/2}$ in all the simulations without loss of generality.

III. RESULTS

A. Vortex formation at weak anisotropy

First, we analyze the effective U(1)-symmetric case in which the sixfold anisotropy of our model is extremely weak. That is, the case in which the phase zero mode is effectively massless and the tangential restoring force in Eq. (9) is

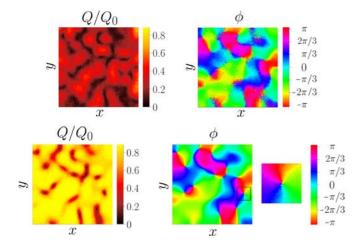


FIG. 2. Snapshots of the amplitude Q (left) and the phase ϕ (right) of the order parameter in the weak-anisotropy case for $\epsilon_f = -0.7$. They illustrate the distribution of vortices at the formation time τ_0 (primordial vortices, top) and at the vortex consolidation time τ_1 (bottom). The zoomed area emphasizes the U(1) character of the vortices in this case.

negligible. The evolution of the vortex pattern obtained in this case is illustrated in Fig. 2.

In our analysis, we first define the time τ_0 at which primordial vortices form as the time at which a pattern of domains of well-defined uniform and stable phase ϕ shows up for the first time. This is the relaxation time of the phase ϕ , and can be determined from the tangential force f_{ϕ} acting on ϕ [see Eq. (9)]. The behavior of the sample-averaged strength of this force $\langle |f_{\phi}| \rangle$ as a function of time is illustrated in Fig. 3(a) for $\epsilon_f =$ -0.3. As we can see, this force drops to a small asymptotic value signaling the formation of metastable domains. In such a quasistationary state, only the stochastic component of Eq. (9) survives. Therefore, since the anisotropic tangential restoring force term in Eq. (9) is negligible, we infer that it is the tangential tension force that causes the relaxation of ϕ until domains of quasiuniform phase get formed. This is nothing but the realization of the geodesic rule in a second-order phase transition, formulated for the first time by Kibble in first-order phase transitions [8,24]. Thus, τ_0 can be identified with the time at which the slope of the tangential force at its saddle point intercepts its long-time asymptote [see Fig. 3(a)]. We use this time to define the density of primordial vortices, $n_0 \equiv n(\tau_0)$, which further evolves in time as illustrated in Fig. 3(b).

To further clarify the evolution of the vortex distribution we track the concomitant changes in the amplitude of order parameter. These changes reveal three dynamical regimes [see Fig. 3(a)]. In a first stage, the linear dynamics of Q is dominated by its diffusion in configuration space as a result of the stochastic force. Diffusion dominates completely the dynamics up to certain *diffusion time*, τ_d , at which Q starts rolling down the effective potential and nonlinear effects become apparent. Diffusion causes an effective delay of the phase transition, whereas the steep descent is caused by the radial restoring force which derives from the effective potential. The phenomenon of *delayed bifurcation* was firstly noticed by Lythe in the dynamics of the phase transition

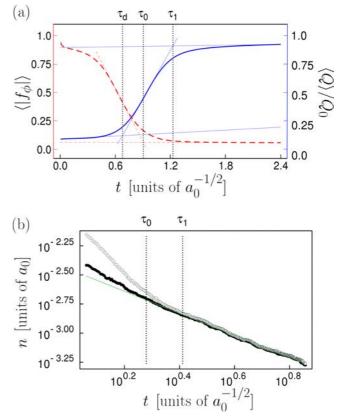


FIG. 3. (a) Sample-averaged tangential force strength (dashed red, in arbitrary units) and normalized amplitude of the order parameter (solid blue) as a function of time, in the weak-anisotropy case, for $\epsilon_f = -0.3$. The vertical dotted lines indicate the diffusion time, τ_d , the phase relaxation time, τ_0 , and the vortex consolidation time, τ_1 , as defined in the main text. (b) Graphical determination of the corresponding vortex density as a function of time (log-log scale). The same power-law behavior is obtained according to two different methods (solid and open circles, respectively; cf. Ref. [34]) beyond τ_1 .

of a field theory [35]. However, in contrast with Lythe's approach and other works [29,30], our consistent treatment of the Langevin term causes the control parameter ϵ (analogous to the bifurcation parameter in Ref. [35]) to take its final value much earlier than τ_d . In fact, once the system is below the transition point, the stochastic Langevin force is the dominant force all the way until the diffusion time τ_d is reached. This prevents the formation of actual vortices during the quench towards the final temperature, and thus any scaling with the quench rate. This condition is in fact an alternative definition for the ultrafast character of the quench. Accordingly, the order parameter starts deviating from zero much later than the moment at which the reduced temperature has reached its final value ϵ_f . Subsequently, the amplitude Q eventually reaches its equilibrium value, Q_0 , at any of the six minima of the effective potential, except at those points where the transition is frustrated by the presence of the topological defects. We identify this event with the vortex consolidation time, τ_1 , which is slightly longer than τ_0 . τ_1 is the relaxation time of the amplitude Q, and thus signals the accomplishment of the phase transition. In fact, during the intermediate regime

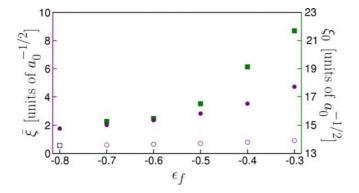


FIG. 4. Average distance between vortices, ξ_0 , as a function of the final temperature of the quench ϵ_f in the weak-anisotropy regime (squares), compared to the mean-field correlation lengths, $\bar{\xi}$ (open circles), and $\bar{\xi}_6$ (solid circles).

between τ_d and τ_1 , Q is mainly driven by the radial restoring force [see Eq. (8)], while in the final stage the stochastic force takes over so that the total radial force f_Q tends to an asymptotic stationary value. The consolidation time τ_1 is thus associated to the crossover between these two regimes [see Fig. 3(a)].

The density of primordial vortices n_0 as a function of the final reduced temperature ϵ_f is summarized in Table I, together with the average distance between those vortices, $\xi_0 \simeq n_0^{-1/2}$, relaxation times, and mean-field values (see Sec. II B). We note that the density of primordial vortices increases monotonically with the decrease of ϵ_f (so that the average distance between vortices decreases). The three characteristic times, in their turn, decrease as the final temperature of the quench does. We interpret that these tendencies have their root in the diffusive dynamics of the order parameter in configuration space, which is enhanced by the temperature. That is, the higher the temperature, the longer the period in which the stochastic force causes Q to fluctuate randomly around Q =0. Hence, τ_d increases with the final temperature. In turn, diffusion slows down the rolling of Q towards any of the minima at $Q = Q_0$, delaying this way the start of the nonlinear dynamics and thus the relaxation processes. The latter implies that τ_0 and τ_1 do also increase with the final temperature. In addition, the persistent thermal fluctuations after the diffusion period do also enhance the periods of relaxation of the phase and the amplitude of the order parameter, causing an increase of the time intervals, $\tau_0 - \tau_d$ and $\tau_1 - \tau_0$, with the final temperature. Altogether, it results in an effective increase of the vortex separation with the final temperature of the quench.

It is also remarkable that the computed vortex correlation lengths, ξ_0 , and relaxation times, τ_0 and τ_1 , differ in an order of magnitude with respect to the values computed in the mean-field linear approximation. Yet, the relative variations of ξ_0 , τ_0 , and τ_1 with the temperature are approximately proportional to the variations of the mean-field correlation lengths and relaxation times, respectively, as illustrated in Fig. 4 for the lengths. The extrapolation of this result to the standard KZ picture implies that the nonuniversal prefactor of the KZ scaling can play a role for the quantitative analysis of the vortex formation. Finally, it is worth noting that, in

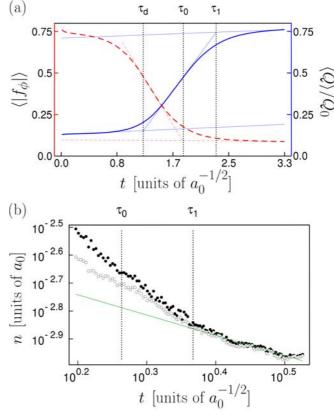


FIG. 5. (a) Sample-averaged tangential force strength (dashed red, in arbitrary units) and normalized amplitude of the order parameter (solid blue) as a function of time in the strong-anisotropy case for $\epsilon_f = -0.3$. The vertical dotted lines indicate the diffusion time, τ_d , the phase relaxation time, τ_0 , and the vortex consolidation time, τ_1 , as defined in the main text. (b) Graphical determination of the corresponding vortex density as a function of time (log-log scale). The same power-law behavior is obtained according to two different methods (solid and open circles, respectively; cf. Ref. [34]) beyond τ_1 .

comparison to previous works in which the amplitude of the noise term is considered small [29,30,35], the vortex density here is not affected by the relationship between that amplitude and the quench rate, but by the relationship of the final temperature and the damping rate.

B. Vortex formation at strong anisotropy

Next, we investigate the impact of the sixfold anisotropy on the formation of vortices. Thus, we consider the extended \mathbb{Z}_6 case described by Eq. (2) in which the phase zero mode becomes massive as described in Secs. II B and II C. As previously mentioned, this is the situation found in hexagonal multiferroic manganites [31–33].

Compared with the previous U(1) situation, the diffusion regime is shortened [see Fig. 5(a)]. This is mainly due to the additional contribution to radial restoring force generated by the anisotropy. However, the phase relaxation interval extends longer due to the anisotropy contribution to the tangential restoring force that opposes to the balance of the tangential tension (i.e., $\tau_0-\tau_d$ increases). As a result, the phase relaxation time τ_0 is approximately equivalent in both cases [see Fig. 5(a)], which further yields a similar density of primordial vortices [see Fig. 5(b)]. Lastly, once the phase gets relaxed, it takes a shorter time for the vortices to consolidate under the action of the additional anisotropic radial restoring force. We note that, due to the increased size of the vortex cores, the sample-averaged amplitude $\langle Q \rangle$ is noticeably smaller than its equilibrium value Q_0 at all times [see Fig. 5(a)]. This was previously pointed out in [33] from an experimental analysis of the static distribution of an order parameter described by the same \mathbb{Z}_6 model.

C. Vortex network evolution

Finally, we analyze the subsequent evolution of the vortex network and, in particular, the vortex-antivortex annihilation process that eventually results in a homogeneous nonsymmetric phase (see Ref. [36] for a movie illustrating the evolution of the complete thermalization process). For this purpose, we fit the time evolution of the vortex density for each of the temperatures of the quench to a power-law function $n(t) \simeq n(\tau_1)/t^{\alpha}$ for $t \ge \tau_1$, as shown in Figs. 3(b) and 5(b). Thus, we find $\alpha \approx 1$ in the U(1) weak-anisotropy case and $\alpha \approx 3/4$ in the \mathbb{Z}_6 strong-anisotropy one. This means that, despite the fact that the overlap between vortices is larger in the strong-anisotropy case for the vortex core is larger, the annihilation rate is slower. This signals the impact of the \mathbb{Z}_6 anisotropy in the short-range vortex-antivortex interaction.

IV. CONCLUSIONS AND DISCUSSION

We have shown that, in ultrafast quenches, the phase transition of a scalar field with a topologically nontrivial vacuum manifold, is accompanied by the formation of a metastable network of topological defects whose density depends on the final temperature of the quench.

The formation process involves three distinguishable and complementary mechanisms. Namely, the diffusive dynamics of the order parameter in configuration space as a result of its coupling to the thermal bath; the local relaxation of its phase as a result of the tension forces between adjacent domains; and the global relaxation of the amplitude of the order parameter as it rolls down the effective potential. Each of these effects possess characteristic times, τ_d , τ_0 , and τ_1 ,

respectively. The primordial vortex network shows up at τ_0 , whereas its consolidation takes place at τ_1 . Hence, the spatial distribution of vortices is not determined by their core radius, but by the correlation length of the phase of the order parameter at τ_0 , ξ_0 . While the alternative choice of τ_1 as the characteristic time for this determination is also possible, it is less physical since then the intertwined dynamics of both the phase and the amplitude have an impact on the resulting topological structure. In fact, from a practical perspective, the unambiguous connection between the phase relaxation time and the density of primordial vortices is one of the main results of our work.

We find that the distance between the primordial vortices increases monotonically with the final temperature of the quench. We explain this behavior as a result of the persistent thermal fluctuations of the order parameter throughout the transition. Moreover, τ_d as well as the intervals between the times τ_d , τ_0 , and τ_1 increase for higher temperatures, which then causes an effective delay of the phase transition together with an increase on the average distance between vortices. In some sense, this can be seen as a sort of ancestor for some intrinsic deviations from the KZ scaling that have been predicted for slow quenches [37].

In the weak-anisotropy regime, once formed, vortices and antivortices annihilate at a rate inversely proportional to time. In turn, all these effects hinder the dissipation of energy, causing a delay in the accomplishment of the thermalization process. Including a sixfold anisotropy we find that, whereas the vortex formation process is unaffected, their annihilation rate slows down, signaling the impact of the \mathbb{Z}_6 anisotropy in the short-range vortex-antivortex interaction.

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