Spontaneous creation of skyrmions in a two-component Bose-Einstein condensate

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We investigate the stability of a vortex ring in a miscible two-component Bose-Einstein condensate confined in a harmonic potential, where the vortex cores in the two components are initially overlapped. Solving the Gross-Pitaevskii equation numerically, we find that the overlapped vortex rings in the two components are dynamically unstable against separation and that they can form linked vortex rings, resulting in a three-dimensional skyrmion. The parameter range for spontaneous skyrmion generation is determined by the Bogoliubov analysis.

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I. INTRODUCTION

Skyrmions are nontrivial topological structures proposed in nuclear physics to explain hadrons [\[1\]](#page-4-0). Skyrmions have since transcended their theoretical origins and have been extended to various physical systems. Numerous theoretical proposals have facilitated the observation and manipulation of skyrmions in chiral magnets [\[2\]](#page-4-0), liquid crystals [\[3\]](#page-4-0), quantum Hall systems [\[4–6\]](#page-4-0), superconductors [\[7,8\]](#page-4-0), semiconductors [\[9\]](#page-4-0), acoustics $[10]$, superfluid ³He $[11,12]$, and Bose-Einstein condensates (BECs) of ultracold gases [\[13,14\]](#page-5-0). The order parameters with internal degrees of freedom enable the atomic BECs to support a wide range of topological entities such as dark solitons $[15]$, quantized vortices $[16,17]$, monopoles [\[18–20\]](#page-5-0), and knots [\[21–23\]](#page-5-0). The system of a multicomponent BEC is thus a suitable platform for examining the behavior of topological spin structures, including skyrmions.

Various studies on skyrmions in multicomponent BECs have been reported. Two-dimensional (2D) skyrmions have been realized experimentally using the Raman process [\[24\]](#page-5-0) and magnetically induced spin rotations [\[25\]](#page-5-0). Threedimensional (3D) skyrmions have been successfully created in a spin-1 BEC using a spin rotation technique [\[26\]](#page-5-0). A wide variety of methods for creating skyrmions in multicomponent BECs have been proposed, including electromagnetically induced transitions [\[13\]](#page-5-0), spin manipulation using a structured magnetic field and laser beams [\[27–30\]](#page-5-0), spin-orbit interaction [\[31,32\]](#page-5-0), annihilation of domain walls [\[33\]](#page-5-0), Plateau-Rayleigh instability [\[34\]](#page-5-0), optical excitation of atoms [\[35\]](#page-5-0), and a moving obstacle [\[36\]](#page-5-0). The stability and dynamics of skyrmions have also been investigated extensively [\[37–49\]](#page-5-0).

In this paper, we show that a 3D skyrmion is generated spontaneously in a two-component BEC. As an initial state, we consider a $U(1)$ vortex ring in a miscible two-component BEC, where quantized vortex rings in the two components are totally overlapped with each other. These overlapped vortex rings are shown to be dynamically unstable against separation. For some parameters, we find that the separated vortices are linked with each other, resulting in a skyrmion with an integer winding number.

The remainder of the paper is organized as follows. A skyrmion in a two-component BEC is introduced in Sec. II. The dynamics of the system and the spontaneous generation of skyrmions are demonstrated in Sec. [III.](#page-1-0) The Bogoliubov analysis to find the parameter range for the skyrmion generation is described in Sec. [IV.](#page-2-0) An experimental scenario to observe the formation of a 3D skyrmion is proposed in Sec. [V,](#page-4-0) and conclusions drawn from this study are offered in Sec. [VI.](#page-4-0)

II. THREE-DIMENSIONAL SKYRMION IN A TWO-COMPONENT BEC

We consider a two-component BEC at zero temperature, represented by the macroscopic wave functions $\Phi_1(\mathbf{r}, t)$ and $\Phi_2(r, t)$ within the mean-field approximation. These twocomponent wave functions can be expressed as

$$
\Psi(r) = \begin{pmatrix} \Phi_1(r) \\ \Phi_2(r) \end{pmatrix} = \sqrt{\rho(r)} \begin{pmatrix} \Lambda_1(r) \\ \Lambda_2(r) \end{pmatrix}, \tag{1}
$$

where $\rho(\mathbf{r}) = |\Phi_1(\mathbf{r})|^2 + |\Phi_2(\mathbf{r})|^2$ is the total density and (Λ_1, Λ_2) satisfies the normalization condition $|\Lambda_1(\mathbf{r})|^2 +$ $|\Lambda_2(\mathbf{r})|^2 = 1$. The two complex numbers $\Lambda_1(\mathbf{r})$ and $\Lambda_2(\mathbf{r})$ represent a pseudospin-1/2 state on the SU(2) manifold. In a two-component BEC, a skyrmion is characterized as a configuration in which the physical space \boldsymbol{r} is continuously and topologically mapped onto the SU(2) manifold. In this mapping, all the points at infinity ($|r| \rightarrow \infty$) are mapped onto the common state (Λ_1, Λ_2) ; therefore, the structure of this mapping is mathematically described by the third homotopy group, $\pi_3[SU(2)] = \mathbb{Z}$ [\[1,](#page-4-0)[14,37,50\]](#page-5-0). This mapping is classified by an integer topological invariant known as the winding number or topological charge, which quantifies its topological nature.

A skyrmion in an equally mixed two-component BEC can be constructed by the following expression,

$$
\Psi(r) = \frac{1}{\sqrt{2}} e^{-if(r)\hat{r}\cdot\sigma} \begin{pmatrix} 1 \\ 1 \end{pmatrix}
$$

=
$$
\frac{1}{\sqrt{2}} \begin{pmatrix} \cos f(r) - i(\cos \theta + \sin \theta e^{-i\phi}) \sin f(r) \\ \cos f(r) + i(\cos \theta - \sin \theta e^{i\phi}) \sin f(r) \end{pmatrix}, (2)
$$

where $f(r)$ is a monotonically decreasing function satisfying the boundary condition $f(0) = \pi$ and $f(\infty) = 0$,

FIG. 1. Skyrmion state defined in Eq. [\(2\)](#page-0-0), which consists of linked vortex rings in two components. (a) Isodensity surface for both components $(|\psi_1|^2 = |\psi_2|^2 = 0.1)$; the color represents the phase at the surface, where the phase changes by 2π around the tubes. (b) and (c) Cross-sectional density (main panels) and phase (insets) profiles for components 1 and 2 on the $y = 0$ plane, respectively.

σ represents the vector of the Pauli matrices, and \hat{r} = $(\sin \theta \cos \phi, \sin \theta \sin \phi, \cos \theta)$ is the unit vector in polar coordinates (r, θ, ϕ) . The rotation matrix in Eq. [\(2\)](#page-0-0) covers whole elements of SU(2); hence it corresponds to a mapping of a 3D unit sphere S^3 parametrized by $[f(r), \theta, \phi]$ onto the $SU(2)$ manifold. Figure 1 shows the skyrmion state defined in Eq. [\(2\)](#page-0-0). This state contains quantized vortex rings in both components, which are linked with each other [\[51\]](#page-5-0), as shown in Fig. 1.

In general, a two-component state can be written as

$$
\Psi(r) = \sqrt{\rho(r)} \begin{pmatrix} e^{i\eta_1(r)} \cos \frac{\kappa(r)}{2} \\ e^{i\eta_2(r)} \sin \frac{\kappa(r)}{2} \end{pmatrix},
$$
(3)

where η and κ are real functions. The winding number *W* for a 3D skyrmion is defined as [\[1,](#page-4-0)[13,14,37,38\]](#page-5-0)

$$
W = \frac{1}{8\pi^2} \int d\mathbf{r} \sum_{\alpha,\beta,\gamma} \epsilon_{\alpha\beta\gamma} \sin \kappa(\mathbf{r}) (\partial_\alpha \kappa) (\partial_\beta \eta_1) (\partial_\gamma \eta_2), \quad (4)
$$

where $\epsilon_{\alpha\beta\gamma}$ represents the antisymmetric tensor and α , β , and γ are summed over *x*, *y*, and *z*. As long as the parameters κ , η_1 , and η_2 are continuous, the winding number *W* is an integer. Equation [\(2\)](#page-0-0) contains a skyrmion structure with $W = 1$.

The winding number is invariant under global spin rotation, which generates topologically equivalent states. Applying spin rotation $e^{i\sigma_y \pi/4}$ to the state in Eq. [\(2\)](#page-0-0), we obtain the skyrmion state discussed in Refs. [\[13,33,36,48\]](#page-5-0), in which component 1 has a vortex ring and its core is occupied by component 2 having 2π rotation of the phase along the ring.

III. NUMERICAL SIMULATIONS OF DYNAMICS

We consider a two-component BEC of a dilute atomic gas with mass *m* trapped inside an external potential *V* (*r*). The dynamics of the system at zero temperature under the mean-field approximation is described by the coupled Gross-Pitaevskii (GP) equations,

$$
i\hbar \frac{\partial \Phi_j}{\partial t} = \left(-\frac{\hbar^2}{2m} \nabla^2 + V(r) + g_{jj} |\Phi_j|^2 + g_{jj'} |\Phi_{j'}|^2 \right) \Phi_j,
$$
\n(5)

where $(j, j') = (1, 2)$ and $(2, 1)$. Here, $\Phi_j(\mathbf{r}, t)$ is normalized as $\int |\Phi_j|^2 d\mathbf{r} = N_j$ with N_j being the number of atoms in the *j*th component. The coefficient $g_{jj'} = 4\pi \hbar^2 a_{jj'}/m$ represents the interaction parameter with $a_{jj'}$ being the *s*-wave scattering length between the *j*th and *j'*th components. In the following, we set the number of atoms in both components to be the same, $N_1 = N_2$. The intercomponent interaction parameter *g*¹² is assumed to be a variable that satisfies miscible condition $g_{11}g_{22} > g_{12}^2$ [\[52\]](#page-6-0). For simplicity, the intracomponent interactions are assumed to be identical: $g_{11} = g_{22} \equiv g > 0$. We also assume that the system is confined in a spherically symmetric trap given by $V(r) = m\omega^2 r^2/2$, where ω is the trap frequency.

We rescale position $\mathbf{r} = \tilde{\mathbf{r}} l_h$, time $t = \omega^{-1} \tau$, and wave function $\Phi_j(\mathbf{r}, t) = \tilde{\Phi}_j(\tilde{\mathbf{r}}, \tau) N_j^{1/2} / l_h^{3/2}$, where $l_h = [\hbar/(m\omega)]^{1/2}$. Equation (5) then takes a nondimensional form,

$$
i\frac{\partial \tilde{\Phi}_j}{\partial \tau} = \left(-\frac{\tilde{\nabla}^2}{2} + \tilde{V} + \tilde{g}|\tilde{\Phi}_j|^2 + \tilde{g}_{jj'}|\tilde{\Phi}_j|^2\right)\tilde{\Phi}_j, \quad (6)
$$

where $\tilde{V} = \frac{\tilde{r}^2}{2}$, $\tilde{g} = \frac{4\pi N_j a}{l_h}$, and $\tilde{g}_{jj'} = \frac{4\pi N_j a_{jj'}}{l_h}$ are the scaled potential and scaled interaction coefficients, respectively. In the following discussion, we omit the tildes from the nondimensional quantities.

We use the split-operator pseudospectral method [\[53\]](#page-6-0) to integrate the 3D GP equation numerically to obtain imaginaryand real-time evolutions. The numerical mesh size is set to $(256)^3$ with a spatial step size of $dx = dy = dz = 0.075$, and the time step is typically $dt = 0.001$. The numerical box is sufficiently larger than the condensate, and the periodic boundary condition imposed by the spectral method does not affect the results.

We prepare the stationary vortex-ring state [\[54–56\]](#page-6-0) as an initial state in which both components have the same wave function containing a quantized vortex ring. To generate this state numerically, we first prepare the ground state without vortices by imaginary-time evolution. We next imprint a circular vortex ring in both components by multiplying the wave functions by $e^{i\Theta(r)}$, where

$$
\Theta(r) = \tan^{-1} \frac{z}{r_{\perp} - R_r} - \tan^{-1} \frac{z}{r_{\perp} + R_r}.
$$
 (7)

Here, $r_{\perp} = \sqrt{x^2 + y^2}$ and R_r is the radius of the vortex ring to be imprinted in both components. We perform a short imaginary-time propagation (duration of 0.3) after the phase $\Theta(r)$ is imprinted on the wave functions to remove the excess energy of the imprinted vortices. If the value of R_r is appropriately chosen, the imaginary-time evolution almost maintains the radii of the vortex rings and the stationary state is achieved. Figure [2](#page-2-0) shows the stationary state with overlapped vortex rings, where $R_r = 3.5l_h$ is used. In Fig. [2\(b\),](#page-2-0) the two density holes indicate the cross section of the vortex ring with 2π phase winding. The two components have the same wave function and the vortex rings in the two components are totally overlapped with each other, having the same radius. A small numerical perturbation is introduced to each component of the initial state before the real-time evolution to break the exact numerical symmetry and trigger the dynamical instability.

Figure [3](#page-3-0) shows the real-time evolution starting from the stationary vortex-ring state in Fig. [2.](#page-2-0) The overlapped vortex rings in the two components are dynamically unstable against separation. This instability is similar to that in two overlapped

FIG. 2. Initial stationary state with overlapped vortex rings in component 1 and component 2 obtained by imaginary-time evolution of the GP equation with $g = 5000$ and $g_{12} = 0.65g$. The two components have the same density and phase profiles. (a) Isodensity surfaces at one-half the peak density, where the outer spherical surface is made transparent for visibility. (b) Cross-sectional density (main panel) and phase (inset) profiles of both components on the $x = 0$ plane. We set $R_r = 3.5l_h$ for preparing the stationary state. The size of the box in (a) is $(19.2l_h)^3$, with the origin at the center. The field of view in (b) is $19.2l_h \times 19.2l_h$, and the units of density are N_j/l_h^3 .

vortices in two components in 2D systems [\[57,58\]](#page-6-0). For a certain range of interaction parameter *g*12, we find that these coaxial vortex rings separate and become linked with each other, which results in the skyrmion. Figures $3(a)-3(c)$ show the dynamics of spontaneous generation of a skyrmion for $g_{12} = 0.65g$. The two vortex rings already become linked immediately after separation ($\tau \simeq 40$), and the linkage is maintained for a long time (until $\tau \simeq 160$). Figure [3\(d\)](#page-3-0) shows the time evolution of the winding number *W* . Since the density outside the condensate is zero and the phase is ill defined, the integral in Eq. [\(4\)](#page-1-0) is taken only for $r < r_{\text{cutoff}} = 5l_h$. The winding number *W* increases from 0 to \approx 1 when the vortex rings separate and become linked. The fluctuations in the winding number during 20 $\lesssim \tau \lesssim$ 40 are due to the ill-defined phases near the separating vortex cores, and those for $\tau \geq 40$ are due to the cutoff r_{cutoff} . At $\tau \simeq 160$, the link between the vortex rings breaks and the winding number returns to $W \simeq 0$ (data not shown). The skyrmion with $W = 1$ or -1 is obtained randomly depending on the initial random noise. In the case of $g_{12} = 0.75g$, spontaneous generation of a skyrmion with double-winding linkages is obtained at $\tau \simeq 40$, as shown in Fig. [3\(e\),](#page-3-0) which has the winding number $W \simeq -2$. After $\tau \simeq 65$, it reduces to the single-winding linkage [Fig. [3\(f\)\]](#page-3-0), followed by the leapfrog dynamics at $\tau \simeq 90$ [Fig. [3\(g\)\]](#page-3-0). For $g_{12} = 0.60g$, no skyrmion is generated and only the leapfrog dynamics is observed (Fig. [4\)](#page-3-0). The results show that the separated vortex rings are not linked with each other and that the axial symmetry is retained for a long time. The winding number *W* for this case remains zero in the time evolution.

We performed systematic simulations using the GP equation for various values of *g*¹² and found that the skyrmion with $|W| \approx 1$ is generated for $0.62g \lesssim g_{12} \lesssim 0.70g$. The leapfrog dynamics of the vortex rings are obtained for $g_{12} \lesssim 0.61g$. For $0.74g \lesssim g_{12} \lesssim 0.77g$, we obtain the skyrmion with $|W| \simeq 2$ with a short lifetime, followed by the leapfrog dynamics.

IV. BOGOLIUBOV ANALYSIS

We here use the Bogoliubov analysis to investigate the dynamical stability of the system. We separate the wave function as

$$
\Phi_j(\mathbf{r}, \tau) = [\psi_j(r_\perp, z) + \delta \Phi_j(\mathbf{r}, \tau)] e^{-i\mu_j \tau}, \quad (8)
$$

where ψ_j is the stationary state, $\delta \Phi_j$ is the small deviation, and μ_i is the chemical potential. Substituting Eq. (8) into the GP equation [\(6\)](#page-1-0) and neglecting the second- and third-order terms in $\delta \Phi_j$, we obtain ($j \neq j'$)

$$
i\frac{\partial(\delta\Phi_j)}{\partial \tau} = \left(-\frac{\nabla^2}{2} + V - \mu_j\right)\delta\Phi_j
$$

+ $g_{jj}(2|\psi_j|^2\delta\Phi_j + \psi_j^2\delta\Phi_j^*)$
+ $g_{jj'}(|\psi_{j'}|^2\delta\Phi_j + \psi_{j'}^*\psi_j\delta\Phi_{j'} + \psi_{j'}\psi_j\delta\Phi_{j'}^*).$
(9)

We write the deviation $\delta \Phi_j$ as

$$
\delta\Phi_j(\mathbf{r},\tau) = u_j(r_\perp, z)e^{i\ell\phi}e^{-i\lambda\tau} + v_j^*(r_\perp, z)e^{-i\ell\phi}e^{i\lambda^*\tau}, \quad (10)
$$

where an integer ℓ denotes the angular momentum of the excitation. Substituting Eq. (10) into Eq. (9) , we obtain the Bogoliubov–de Gennes equation ($j \neq j'$),

$$
\begin{aligned}\n&\left[-\frac{1}{2}\left(\nabla_{\perp}^{2} + \nabla_{z}^{2} - \frac{\ell^{2}}{r_{\perp}^{2}}\right) + V - \mu_{j}\right]u_{j} + g_{jj}(2|\psi_{j}|^{2}u_{j} + \psi_{j}^{2}v_{j}) \\
&+ g_{jj'}(|\psi_{j'}|^{2}u_{j}\delta\psi_{j} + \psi_{j}^{*}\psi_{j}u_{j'} + \psi_{j'}\psi_{j}v_{j'}) = \lambda u_{j}, \\
&\left[-\frac{1}{2}\left(\nabla_{\perp}^{2} + \nabla_{z}^{2} - \frac{\ell^{2}}{r_{\perp}^{2}}\right) + V - \mu_{j}\right]v_{j} \\
&+ g_{jj}(2|\psi_{j}|^{2}v_{j} + \psi_{j}^{*2}u_{j}) \\
&+ g_{jj'}(|\psi_{j'}|^{2}v_{j} + \psi_{j'}\psi_{j}^{*}v_{j'} + \psi_{j'}^{*}\psi_{j}^{*}u_{j'}) = -\lambda v_{j}.\n\end{aligned} \tag{11b}
$$

Numerically, we obtain the stationary state $\psi_i(r_\perp, z)$ containing vortices using the imaginary-time evolution of the GP equation, in a manner similar to Fig. 2, where r_{\perp} and *z* are discretized with the step size $dr_{\perp} = dz = 0.075$. We next refine the stationary state using the Newton-Raphson method [Fig. $5(a)$]. Using this stationary state, we calculate the eigenvalues and eigenfunctions of Eq. (11) using the Lanczos method. The existence of at least one eigensolution with a positive imaginary part (Im $\lambda > 0$) indicates dynamic instability because the corresponding excitation mode exhibits exponential growth in time.

An example of the eigenfunctions u_j and v_j for a dynamically unstable mode is shown in Figs. $5(b)$ – $5(e)$. Since Eq. (11) does not depend on the sign of ℓ , we show only the results for $\ell \geqslant 0$. The eigenfunctions u_j are localized near the vortex core [Figs. $5(b)$ and $5(c)$], whereas v_j are distributed [Figs. [5\(d\)](#page-3-0) and [5\(e\)\]](#page-3-0). They satisfy $\int dr \sum_{j=1}^{2} (|u_j|^2 - |v_j|^2) =$ 0. We note that the localized functions satisfy $u_1 = -u_2$, which leads to vortex separation, as discussed below. The eigenfunctions for other unstable modes have similar shapes. There is also a dynamically unstable mode in which u_j and v_j are exchanged, since (λ, u_j, v_j) and $(-\lambda^*, v_j^*, u_j^*)$ are conjugate solutions.

FIG. 3. Dynamics of spontaneous skyrmion generation obtained by the GP equation for $g = 5000$, where the initial state is the overlapped vortex rings (Fig. [2\)](#page-2-0). Isodensity surfaces for component 1 (yellow or light gray) and component 2 (blue or dark gray) for (a)–(c) $g_{12} = 0.65g$ and (e)–(g) $g_{12} = 0.75g$. See the Supplemental Material for animations of the dynamics in (a)–(c) and (e)–(g) [\[59\]](#page-6-0). (d) and (h) Time evolution of the winding number *W* for $g_{12} = 0.65g$ and $g_{12} = 0.75g$, respectively.

The integer ℓ for the unstable mode corresponds to the winding number *W* for the generated state, which can be understood as follows. Suppose that the core of the vortex ring is located at $r_{\perp} = r_0$ and $z = 0$, as shown in Fig. 5(a). We focus on the region very close to the core of the vortex ring, where u_i are dominant in the excitation. Approximating that u_i are uniform near the vortex core, we can express the time evolution of the wave function as

$$
\psi_j \propto r_\perp - r_0 + iz + C_j(t)e^{i\ell\phi},\tag{12}
$$

where C_i are proportional to the values of u_i at the vortex core. Without loss of generality, C_i can be taken to be real, and Eq. (12) becomes

$$
\psi_j \propto r_\perp - (r_0 - C_j \cos \ell \phi) + i(z + C_j \sin \ell \phi). \tag{13}
$$

Thus, the position of the vortex core is displaced by $(-C_j \cos \ell \phi, -C_j \sin \ell \phi)$; i.e., the displaced vortex ring winds ℓ times around the original ring from $\phi = 0$ to 2π , which corresponds to the winding number $W = \ell$ for the skyrmion. Since $u_1 = -u_2$ (i.e., $C_1 = -C_2$), the vortices in the two components separate from each other.

FIG. 4. Leapfrog dynamics obtained by solving the GP equation for $g = 5000$ and $g_{12} = 0.60g$. (a) $\tau = 45$, (b) $\tau = 60$, and (c) $\tau = 90$. Isodensity surfaces for both components are shown. See the Supplemental Material for an animation of the dynamics [\[59\]](#page-6-0).

Figure 5(f) shows the imaginary part of the excitation frequency, Im λ , for $\ell = 0$, 1, and 2 as a function of g_{12}/g . The system is dynamically unstable in the entire range shown, $0.55 \lesssim g_{12}/g \lesssim 0.9$, since there are excitation frequencies

FIG. 5. Bogoliubov analysis for $g = 5000$. (a) Density profile for the stationary state ψ_i . (b)–(e) Density profiles for dynamically unstable modes u_1 , u_2 , v_1 , and v_2 , respectively, for $g_{12} = 0.65g$ and $\ell = 1$. (f) Imaginary part of the Bogoliubov excitation frequency Im λ for eigenmodes with $\ell = 0, 1$, and 2 plotted as a function of *g*₁₂/*g*. The vertical dotted lines indicate $g_{12}/g = 0.60$, $g_{12}/g = 0.65$, and $g_{12}/g = 0.75$, corresponding to the parameters used in Figs. 3 and 4.

with a positive imaginary part for any values of *g*12/*g*. For $g_{12}/g \simeq 0.6, 0.65,$ and 0.75, the excitation frequency λ has the largest imaginary part for $\ell = 0, 1$, and 2 [see the plots at the vertical dotted lines in Fig. $5(f)$]. This result is consistent with the dynamics in Figs. [3](#page-3-0) and [4,](#page-3-0) where the generated states have winding numbers $|W| = 0$, 1, and 2 for $g_{12}/g \approx 0.6$, 0.65, and 0.75, respectively. Although Im λ also becomes nonzero for the modes of $\ell \geq 3$, the values of Im λ for $\ell \geq 3$ are always smaller than those for $\ell \leq 2$. We have confirmed that the results are almost unchanged when the spatial step size is doubled, which indicates that the numerical error due to the spatial resolution is unimportant.

V. EXPERIMENTAL PROPOSAL

We consider a two-component BEC of $87Rb$ atoms confined in an isotropic harmonic potential $V = m\omega^2(x^2 + y^2 +$ z^2)/2 with a trap frequency $\omega = 2\pi \times 100$ Hz. In this case, the units of length and time are $l_h = 1.08 \mu m$ and $\omega^{-1} = 1.59$ ms, respectively. In Secs. [III](#page-1-0) and [IV,](#page-2-0) for simplicity, we assumed $g_{11} = g_{22}$, which can be realized by, e.g., using the hyperfine states $|F, m_F\rangle = |1, 1\rangle$ and $|1, -1\rangle$ for the two components. The *s*-wave scattering lengths for these states are $a_{11} = a_{22} = 100.4a_0$ [\[60,61\]](#page-6-0), where a_0 is the Bohr radius. The interaction parameter $g = 5000$ corresponds to $N_1 = N_2 \simeq$ 8×10^{4} .

The initial state in which the two components contain the overlapped vortex rings (Fig. [2\)](#page-2-0) can be prepared as follows. First, a single-component BEC is created and a stationary vortex ring is produced in the BEC. The vortex ring can be created using the phase imprinting technique, which gives a spatially dependent phase to the wave function by laser beams or a magnetic field [\[13](#page-5-0)[,62\]](#page-6-0). A stirring external potential can also produce a vortex ring in the BEC [\[63\]](#page-6-0). Next, one-half of one component is transferred to the other component through the Rabi oscillation using a $\pi/2$ pulse, which results in the overlapped vortex rings in the two-component BEC.

We have numerically confirmed through the GP simulations that similar results can be obtained for the hyperfine states $|1, 1\rangle$ and $|1, 0\rangle$ of ⁸⁷Rb atoms, where the intracomponent scattering lengths $a_{11} = 100.4a_0$ and $a_{22} = 100.86a_0$ are different. In the numerical simulation, we prepared a stationary vortex-ring state for a BEC of the $|1, 1\rangle$ state and then one-half of the atoms were transferred to the $|1, 0\rangle$ state. We then carried out the real-time evolution in a manner similar to Fig. [3.](#page-3-0) We found that the skyrmion is generated for $0.77g_{11} \lesssim$ $g_{12} \lesssim 0.79 g_{11}.$

VI. CONCLUSIONS

We have shown that a 3D skyrmion is spontaneously generated in a two-component miscible BEC trapped in a harmonic potential, in which the intercomponent interaction *g*¹² is controlled. For a certain range of the intercomponent interaction parameter, the initially overlapped vortex rings exhibit dynamical instability and separate into two linked vortex rings (Fig. [3\)](#page-3-0). We carried out the Bogoliubov analysis and numerically obtained the dynamically unstable spectrum (Fig. [5\)](#page-3-0). This phenomenon can be realized in a realistic experimental system using, e.g., a two-component BEC with a feasible number of atoms (\sim 10⁵ atoms of ⁸⁷Rb).

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