Density Dependent Gauge Field Inducing Emergent Su-Schrieffer-Heeger Physics, Solitons, and Condensates in a Discrete Nonlinear Schrödinger Equation

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We investigate a discrete nonlinear Schrödinger equation with dynamical, density-difference-dependent gauge fields. We find a ground-state transition from a plane wave condensate to a localized soliton state as the gauge coupling is varied. Interestingly we find a regime in which the condensate and soliton are both stable. We identify an emergent chiral symmetry, which leads to the existence of a symmetry-protected zero-energy edge mode. The emergent chiral symmetry relates low and high energy solitons. These states indicate that the interaction acts both repulsively and attractively.

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Introduction.—The nonlinear Schrödinger equation (NLSE) is an essential model for a variety of physical systems [1,2]. In ultracold atomic physics, the NLSE describes Bose-Einstein condensates (BECs) in the dilute limit where only the leading order interaction is relevant [3–6]. In optical systems, the NLSE describes light propagating in nonlinear media, originally employed to describe self-focusing beams of light [7–9]. Since these initial investigations a wide range of studies into the NLSE and its extensions have been developed, identifying a variety of nonlinear phenomena [10–16]. Additionally, extensions have been made to simulate gauge theories [17–21]. Much work has gone into understanding the stability of NLSE solutions. This is in part because the onset of instability can often be associated with a phase transition to a stable nondispersive bound state, termed a soliton [22–27].

Solitons and other localized states have been predicted and verified in the NLSE and its extensions [28–31]. These bound states result from the competition between the nonlinearity and the kinetic dispersion. In the case of the discrete NLSEs [32,33], the existence of bound states termed discrete breathers has been shown to be generic [34]. Such self-stabilizing solutions are widely studied for both their mathematical properties as well as their potential applications in developing communications technology. Recently, matter wave breathers were experimentally observed by quenching the strength of the attractive interactions in a quasi-one-dimensional Bose-Einstein condensate [35]. Engineering and manipulation of soliton modes has been achieved through various methods, including time-dependent modulation and spatial modulation [36-40].

An interesting avenue explored in recent years has been to consider NLSE which contains a *dynamical*, *density-dependent*, *gauge field*. Here, in the context of discrete (lattice) models, the gauge field is a nontrivial phase factor one has upon hopping between adjacent sites, and densitydependent gauge field means that the phase factor depends on the density of sites involved in the hopping [41–46]. It has been shown that density-dependent hopping induced by sinusoidal temporal modulations of the nonlinearity in the discrete NLSE leads to the formation of compactons, i.e., nonlinear bound states that, unlike breathers, do not have exponential decay tails [31]. Recently, two of us proposed a novel type of density-dependent gauge field, in the context of interacting non-Hermitian quantum physics, where the hopping depends on the difference of the density of sites involved in hopping [47]; such a gauge field has a unique feature that the sign of the interaction is not determined a priori but dynamically depends on the density gradient of the wave function profile. In this paper, we find that the NLSE derived from such a density-difference-dependent gauge field results in a variety of unique phenomena, including the coexistence of plane-wave condensate state and soliton state, the coexistence of attractive and repulsive solitons, and the emergent Su-Schrieffer-Heeger (SSH)-like physics with a symmetry-protected zero-energy mode. Our paper opens an avenue to explore nonlinear physics of this exotic form of gauge field.

The discrete density difference dependent nonlinear Schrodinger equation (D⁴ NLSE).—We consider a one-dimensional discrete NLSE derived from the following classical Hamiltonian:

$$H = \sum_{j} \Psi_{j+1}^* \left[-J + \gamma (n_{j+1} - n_j) \right] \Psi_j + \text{c.c.}, \quad (1)$$

where Ψ_j is the complex-valued amplitude of the wave function at site j, $n_j = \Psi_j^* \Psi_j$ is the corresponding density,

J is the density-independent part of hopping between adjacent sites, γ is the coupling constant of the density-difference-dependent hopping which can generally be a complex number. We assume that the wave function is normalized as $\sum_j |\Psi_j|^2 = 1$. A crucial aspect of this Hamiltonian is that the sign of the interaction, either repulsive or attractive, is not determined *a priori*, but rather determined dynamically through the density profile of the wave function.

The corresponding NLSE can be obtained by computing the Poisson bracket $i(d\Psi_i/dt) = {\Psi_i, H} = (\delta H/\delta \Psi_i^*)$:

$$i\frac{d\Psi_{j}}{dt} = \{-J + \gamma(2n_{j} - n_{j-1})\}\Psi_{j-1} - \gamma\Psi_{j}^{2}\Psi_{j+1}^{*} + \{-J - \gamma^{*}(2n_{j} - n_{j+1})\}\Psi_{j+1} + \gamma^{*}\Psi_{j}^{2}\Psi_{j-1}^{*}, \quad (2)$$

which is the one-dimensional discrete density difference dependent (D⁴) nonlinear Schrödinger equation we are going to explore in this paper.

We note that the Hamiltonian in Eq. (1) can also be thought of as quantum Hamiltonian by identifying Ψ_j^* and Ψ_j as creation and annihilation operators of a particle at site j. Then the Poisson bracket mentioned above should be replaced by a commutator, and D^4 NLSE describes a situation where the mean-field theory is applicable and the Gross-Pitaevskii treatment of the condensate wave function is justified.

Ground-state phase diagram.—We explore the stationary state of the D⁴ NLSE evolving in time as $\Psi_i(t)$ = $e^{-i\mu t}\Psi_i(0)$, where μ is the chemical potential. Of particular interest is the ground state, which is the state with the lowest energy. We can numerically explore the ground state by the method of imaginary-time propagation, which is to start from a random initial state and simulate the time evolution of the D⁴ NLSE with an imaginary time, that is, to consider the equation obtained after setting $t = -i\tau$ and simulate the evolution with respect to τ . After evolving for a long enough time in τ , one converges to the ground state (if the ground state is unique) [48,49]. Assuming periodic boundary condition and applying an imaginary time evolution method to various values of γ , we see that the ground state is either a plane-wave condensate state or a sharply localized solitonlike state, which we explain now.

We first note that, for states with a uniform density, $n_j = n_{j+1}$, γ dependence in the Hamiltonian vanishes. The plane-wave state e^{ikx} with $k = 2\pi(\text{integer})/L$, where L is the number of lattice sites, is thus a valid stationary state with chemical potential $\mu = -2J\cos(k)$. Among the plane-wave solutions, the ground state is the k = 0 state, and the corresponding chemical potential as well as the energy is $\mu = E = -2J$.

Now the question to ask is if one can obtain a state with lower energy by allowing the density to vary. We indeed find that the system can host localized stationary states which we call solitons. Essential features of the soliton can be captured through the following ansatz where the soliton is spread over only three sites around a site n:

$$\Psi_{n\pm 1} = e^{i\phi_{\pm}} \sqrt{(1-\alpha)/2}, \qquad \Psi_n = \sqrt{\alpha}, \qquad (3)$$

where ϕ_-, ϕ_+, α are variational parameters which we take as real numbers. We expect that the interaction energy dominates in such soliton states. The transition between the k = 0 plane wave and the localized ansatz can be solved analytically for purely real γ and is predicted to occur at $|\gamma| = 2.53J$. For complex γ , the critical $|\gamma|$ weakly depends on its phase, decreasing slightly as the phase goes towards purely imaginary γ where the transition occurs around $|\gamma| \approx 2.5J$. We find from imaginary time evolution that the transition occurs at $|\gamma| \approx 2.04J$ for fully real γ and $|\gamma| \approx$ 1.9 for fully imaginary γ ; these values are smaller than the prediction from the three-site ansatz we obtained above because the soliton state around the transition point can be more spread and/or asymmetric. We numerically confirm that, typically, more than 95% of the weight of the soliton lies on three sites, and the three-site ansatz becomes more and more accurate as $|\gamma|$ becomes larger.

We provide a more detailed analysis for the case of purely imaginary $\gamma = i\gamma_I$. The results are qualitatively the same for any fixed phase of γ with slight quantitative differences as mentioned above. In Fig. 1, we plot the energy and the chemical potentials for the lowest and highest energy localized soliton states as a function of $|\gamma_I|$ obtained from imaginary time evolution. These results were

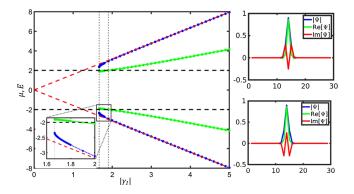


FIG. 1. Left: energy (green solid curve) and the chemical potential (blue dotted curve) of the localized soliton state, in units of J, obtained from the imaginary time propagation method, as a function of the gauge coupling which we assume to be purely imaginary, $\gamma = i\gamma_I$. The red dashed lines correspond to the chemical potential of the localized states obtained when setting J=0. The dashed horizontal black lines correspond to the lowest and highest plane-wave condensate states at $\mu=E=\pm 2J$. The inset highlights the transition point and soliton disappearance. Right: ground-state and highest energy state wave function for J=1, $\gamma=-4.0i$ and length L=29. The blue, green, and red curves are the wave function amplitude, real part, and imaginary part, respectively.

found in excellent agreement also with the ones obtained by means of self-consistent numerical diagonalizations of the nonlinear eigenvalue problem associated with Eq. (2) [30]. The symmetry around E = 0 is a consequence of an emergent chiral symmetry which we discuss later. At $|\gamma_I| > 1.9J$, the solitons are the lowest energy states. In the region $1.67J < |\gamma_I| < 1.9J$, the soliton solution still exists as a stationary state, but its energy is higher than the energy of the plane-wave condensate state. As $|\gamma_I|$ approaches $\approx 1.67J$ from above, the chemical potential of the soliton changes rapidly, and at $|\gamma_I| < 1.67J$, the soliton state no longer exists as a stationary state but rather merges with the extended modes. The energy and the chemical potential of the soliton state just above $|\gamma_I| \approx$ 1.67*J* are $E \approx -1.9J$ and $\mu \approx -2.3J$. Note that we have studied this model in the continuum and found that in solving for a localized state, the energy is unbounded from below, suggesting that soliton formation is a genuine lattice effect (see the Supplemental Material [50]).

Stability of the condensate.—We have just seen that the ground state changes from the plane-wave condensate state to the localized soliton state as one increases the coupling $|\gamma|$. The mechanism behind the formation of the soliton ground state is qualitatively different from more conventional soliton formation in, for example, NLSE with a cubic nonlinearity, where the development of modulational instability of the plane-wave condensate state leads to the formation of solitons [51]. In contrast, in the D⁴ NLSE, there can be regions in the parameter space where the zero momentum plane-wave condensate and the soliton both exist as stable stationary states. Depending on the value of γ , the condensate can become unstable, as we show soon, but the instability of the condensate and the formation of

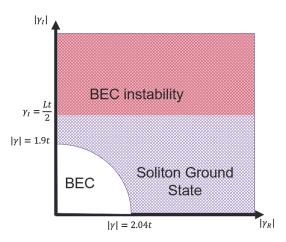


FIG. 2. Phase diagram of the D³NLSE as a function of the complex gauge field coupling γ . The complex gauge coupling is defined as $\gamma_R + i\gamma_I$, J is the single particle hopping parameter, and L is the length of the chain. Note that this phase diagram only considers the linear stability. Numerically we have observed that the instability can occur for completely real γ , likely due to higher order instabilities.

the soliton appear to be not directly related. We now explore this linear stability of the condensate phase and present a complete phase diagram, shown in Fig. 2, including the stability analysis. (In the Supplemental Material [50], we also present the continuum model and demonstrate that its plane-wave solutions have the same stability condition).

To investigate the linear stability of a condensate upon small perturbation, we introduce the Fourier transformed Hamiltonian. Defining the wave function in momentum space with momentum k by $\tilde{\Psi}_k$ as

$$\Psi_j = \frac{1}{\sqrt{L}} \sum_k \tilde{\Psi}_k e^{-ijk},\tag{4}$$

the Hamiltonian is

$$H = \sum_{k} -2J\cos k\tilde{\Psi}_{k}^{*}\tilde{\Psi}_{k}$$

$$+ \frac{1}{L} \sum_{k,k',q} V(q,k)\tilde{\Psi}_{k}^{*}\tilde{\Psi}_{k'}^{*}\tilde{\Psi}_{k+k'-q}\tilde{\Psi}_{q}$$

$$V(q,k) = 2i\gamma_{R}(\sin k - \sin q) + 2i\gamma_{I}(\cos q - \cos k), \quad (5)$$

where we have introduced the notation $\gamma = \gamma_R + i\gamma_I$. The time evolution equation in momentum space is then

$$i\dot{\tilde{\Psi}}_{k} = -2J(\cos k - 1)\tilde{\Psi}_{k} + \sum_{q,k'} V(q,k)\tilde{\Psi}_{k'}^{*}\tilde{\Psi}_{k+k'-q}\tilde{\Psi}_{q}$$
$$+ \sum_{q,k'} V(q,k')\tilde{\Psi}_{k'}^{*}\tilde{\Psi}_{k+k'-q}\tilde{\Psi}_{q}. \tag{6}$$

We assume that the condensate takes place at zero momentum, and add a small perturbation at momentum p to investigate its stability; we set $\tilde{\Psi}_0 \approx 1$, $\tilde{\Psi}_{\pm p} = \delta \tilde{\Psi}_{\pm p}$ which should be kept up to linear order, and $\tilde{\Psi}_k = 0$ otherwise. Inserting them into Eq. (6) and ignoring terms of higher order in $\delta \tilde{\Psi}_{\pm p}$, we obtain the equation for the time derivative of $\delta \tilde{\Psi}_p$ as

$$i\frac{d\delta\tilde{\Psi}_{p}}{dt} = -2J(\cos p - 1)\delta\tilde{\Psi}_{p} + \frac{4i\gamma_{I}}{L}(\cos p - 1)\delta\tilde{\Psi}_{-p}^{*}, \quad (7)$$

and a similar equation for the time evolution of $\delta \tilde{\Psi}_{-p}$. This set of equations can be diagonalized, in a manner analogous to the Bogoliubov transformation, to obtain the dispersion relation, which is

$$\epsilon(p) = 2J\sqrt{\left(1 - \frac{4\gamma_I^2}{J^2 L^2}\right)}(1 - \cos p). \tag{8}$$

This dispersion relation contains unusual features. First, at small momentum p, the dispersion is *quadratic* rather than

a more standard linear excitation spectrum in NLSE. The effect of the interaction γ is to modify the overall multiplicative factor, which is analogous to modifying the effective mass of a particle. The second point to notice is that only the imaginary part of γ enters the dispersion relation. The real part does not affect the excitation properties of the zero momentum condensate phase. The third point is that the dispersion becomes imaginary, implying the appearance of modulational instability, above a threshold value of $\gamma_I > JL/2$; notably the transition depends on the length of the system L. The threshold to the instability of this condensate is drawn in the phase diagram (Fig. 2). We point out that the stability condition is solely determined by γ_I whereas the transition from the condensate ground state to the soliton ground state depends on $|\gamma|$, implying that these two transitions occur from different mechanisms. (The stability condition for a condensate at arbitrary momentum k is presented in the Supplemental Material [50] derived from the Bogoliubov analysis of a bosonic system.)

Chiral symmetry and zero mode.—In Fig. 1, we see that the figure is symmetric around E = 0. This symmetry is a consequence of the chiral symmetry present in the system. By examining the equation of motion, we see that for a stationary state $\{\Psi_i\}$, we obtain a state with the opposite chemical potential by multiplying the state by $e^{i\pi j}$, i.e., a factor of -1 to every other site. In Fig. 1, the wave functions of the ground state and the highest energy state are plotted for $\gamma = -4i$ to demonstrate that they are precisely related by the transformation $e^{i\pi j}$. Note that the existence of these two bound states is a signature of the attractive and repulsive nature of our interaction. Typically the appearance of low energy bound states is a signature of an attractive interaction [25,52] whereas the existence of high energy bound states is a signature of a repulsive interaction [53].

The presence of such a chiral symmetry suggests the existence of a symmetry-protected zero mode analogous to the SSH model. The SSH model is the one-dimensional tight-binding model with alternating weak and strong hopping strengths, and its momentum-space topology is nontrivial when the hopping from the edge starts with a weak hopping. The topologically nontrivial SSH model hosts a zero-energy edge state protected by the chiral symmetry whose wave function takes nonzero values in every other site. Our D⁴ NLSE does not have a priori alternating hopping; however, since the density dependence affects the hopping strength, one can expect an emergent SSH-like physics induced dynamically by an appropriate density distribution. To explore such emergent SSH-like physics, we look for a zero-energy edge mode. Such a zero-energy edge mode must have nonzero wave function in every other site. We consider an open boundary condition with $j \ge 1$, and set the wave function on sites with even indices to zero $\Psi_{2n} = 0$. The D⁴ NLSE

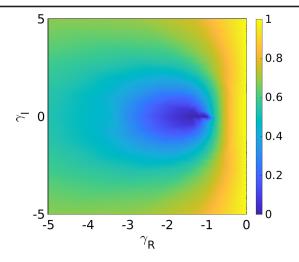


FIG. 3. Localization ratio $|\Psi_3|^2/|\Psi_1|^2$ in the complex γ plane. The coloring corresponds to this ratio. Note that the ratio vanishes at $\gamma = -1$ where the singleton state stabilizes.

taking zero chemical potential leads to the following condition to be satisfied by the wave functions in odd sites:

$$(J + \gamma |\Psi_{2n-1}|^2)\Psi_{2n-1} = (-J + \gamma^* |\Psi_{2n+1}|^2)\Psi_{2n+1}.$$
 (9)

One can iteratively solve this equation for decreasing solution starting from some value of Ψ_1 for $\gamma_R < 0$ or Ψ_L for $\gamma_R > 0$. When the total number of sites is odd, the iteratively constructed solution is the exact zero-energy solution of the D⁴ NLSE. When the total number of sites is even, this construction fails as the final even site with only one adjacent odd site cannot properly satisfy the above equation. We note that there is a zero-energy mode completely localized at one site on the edge, i.e., n=1, when $\gamma/J=-1$, which we call a singleton state. (This state exists regardless of the number of sites). As γ moves from $\gamma/J = -1$, the state becomes more and more delocalized as shown in Fig. 3, which plots the ratio of the $|\Psi_3|^2/|\Psi_1|^2$ in the complex γ plane as a measure of localization. The predominant trends show that as $|\gamma|$ moves from 1 and as γ_I grows the state spreads further. In fact for a purely imaginary γ , the state is evenly spread on odd numbered sites.

Floquet realization.—Finally, we present one method using Floquet engineering to realize the D⁴ NLSE, which can be implemented in systems such as Bose-Einstein condensates of ultracold quantum gases and waveguide arrays. We derive an effective Hamiltonian of a three step Floquet protocol as derived in the formalism presented in Ref. [54]. We consider the time dependent quantum Hamiltonian with periodic drive $V(t) = V(t + T_{\omega})$

$$H(t) = -J \sum_{j} \hat{\Psi}_{j+1}^{\dagger} \hat{\Psi}_{j} + \hat{\Psi}_{j}^{\dagger} \hat{\Psi}_{j+1} + V(t)$$
 (10)

$$V(t) = \begin{cases} V_1, & 0 \le t < \frac{T_{\omega}}{3} \\ V_2, & \frac{T_{\omega}}{3} \le t < \frac{2T_{\omega}}{3} \\ -V_1 - V_2, & \frac{2T_{\omega}}{3} \le t < T_{\omega}, \end{cases}$$
(11)

which corresponds to a free boson with pulsed modulation of the hopping parameter $V_1 = \sum_j J_1 \hat{\Psi}_{j+1}^\dagger \hat{\Psi}_j + \text{H.c.}$ and a pulsed on site interaction $V_2 = U \sum_j \hat{\Psi}_j^\dagger \hat{\Psi}_j^\dagger \hat{\Psi}_j^\dagger \hat{\Psi}_j \hat{\Psi}_j$ in three steps of equal duration with total period $T_\omega = 2\pi/\omega$. In the limit of large ω (i.e., U/ω , $J_1/\omega \ll 1$), the effective Hamiltonian to first order in $1/\omega$ is given as

$$H_{\text{eff}} = \sum_{j} \hat{\Psi}_{j+1}^{\dagger} \left[-J + \frac{2i\pi U J_{1}}{27\omega} (\hat{n}_{j} - \hat{n}_{j+1}) \right] \hat{\Psi}_{j} + \hat{\Psi}_{j}^{\dagger} \left[-J - \frac{2i\pi U J_{1}^{*}}{27\omega} (\hat{n}_{j} - \hat{n}_{j+1}) \right] \hat{\Psi}_{j+1},$$
(12)

which maps directly onto Eq. (1) with $\gamma = 2i\pi U J_1/(27\omega)$ and identifying the quantum operators by classical variables assuming mean-field theory. We control the phase of the complex coupling γ by changing the phase of the modulated hopping J_1 . Complex hopping of J_1 can be engineered through modulation such as in Struck *et al.* [55], which can be combined with our proposed Floquet protocol, or through Raman-assisted tunneling [56]. Similar Floquet interactions have been previously studied in various ultracold and optical systems [41,43,44,57].

Conclusion.—In this Letter, we have presented the D⁴NLSE and its phase diagram. We have derived the conditions for plane-wave stability as well as a phase transition from the extended k=0 plane-wave ground state to a localized ground state. Contrary to conventional NLSE, this model admits the coexistence of a stable k=0 plane wave and a soliton ground state. We have further demonstrated that the interaction of the D⁴ NLSE acts either attractively or repulsively depending on the state in question. Although the D⁴ NLSE itself is fully translationally invariant, the density distribution results in emergent SSH-like physics, and as a consequence, a zero-energy edge-localized mode exists. Finally we derived a possible method of Floquet engineering to realize the D⁴ NLSE.

Many of the characteristic features of D⁴ NLSE we found in this Letter stem from the fact that the sign of the interaction is not determined *a priori*, but rather dynamically determined by the density distribution of the nonlinear wave. The emergent chiral symmetry analogous to the one discussed in this Letter can be a versatile mechanism to obtain interaction-induced topology.

The dynamical dependence of the sign of the interaction represents a new emerging paradigm in the field of nonlinear physics opening up unexpected and unexplored scenarios at both the classical and quantum levels. In this Letter we have focused on properties of the classical ground state of the mean-field problem under this exotic form of dynamical gauge field. A ground-state phase diagram of the corresponding quantum many-body problems as well as its non-Hermitian generalizations are left for future studies.

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