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Synthetic spin–orbit coupling and topological polaritons in Janneys–Cummings lattices

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The interaction between a photon and a qubit in the Janneys–Cummings (JC) model generates a kind of quasiparticle called polariton. While they are widely used in quantum optics, difficulties in engineering-controllable coupling of them severely limit their applications to simulate spinful quantum systems. Here we show that, in the superconducting quantum circuit context, polariton states in the single-excitation manifold of a JC lattice can be used to simulate a spin-1/2 system, based on which tunable synthetic spin–orbit coupling and novel topological polaritons can be generated and explored. The lattice is formed by a sequence of coupled transmission line resonators, each of which is connected to a transmon qubit. Synthetic spin–orbit coupling and the effective Zeeman field of the polariton can both be tuned by modulating the coupling strength between neighboring resonators, allowing for the realization of a large variety of polaritonic topological semimetal bands. Methods for detecting the polaritonic topological edge states and topological invariants are also proposed. Therefore, our work suggests that the JC lattice is a versatile platform for exploring spinful topological states of matter, which may inspire developments of topologically protected quantum optical and information-processing devices.

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INTRODUCTION

The Janneys–Cummings (JC) model proposed in 1963¹ is a seminal theoretical model treating light–matter interaction with full quantum theory, i.e., the interaction of a quantized electromagnetic field with a two-level atom. This model has been widely applied to many quantum platforms for studying the interaction of a quantized bosonic field with a qubit, which now has become the cornerstone in quantum optics and quantum computation.^{2–7} Furthermore, an interconnected array of multiple JC systems can form a JC lattice,^{8–11} which provides an innovative quantum optical platform for studying condensed matter physics. This is highlighted by previous works which show that coupled JC systems can be used to realize the Bose–Hubbard model and investigate superfluid-to-Mott-insulator phase transition.^{12–14} However, spinful lattices have not been simulated in this platform due to the difficulty in engineering a tunable coupling between different cavities.

On the other hand, the search for topological states of matter in artificial systems recently has become a rapidly growing field of research.^{15–24} Topological states are characterized by topological invariants which are robust to the smooth changes in system parameters and disorders, where topological edge states can be employed for robust quantum transport.^{25,26} Therefore, they hold tremendous promise for fundamental new states of matter as well as for dissipationless quantum transport devices and topological quantum computation.²⁷ One of the key ingredients for generating such states is to realize tunable spin–orbit coupling (SOC). Significant theoretical and experimental progress on realizing synthetic SOC recently have been achieved in ultracold atom systems.^{28–30} This progress stimulates great research interests to explore topological states with ultracold atoms trapped in optical

lattices.^{31–34} However, the limited trapping time and the site-addressing difficulty increase the experimental complexity, i.e., it is generally difficult to have modulable coupling between two neighboring sites in an optical lattice.

Here, we find that the JC lattice system can be used to realize various topological spin–lattice models, where synthetic polaritonic SOC and Zeeman field can be induced with in situ tunability, which provides a flexible platform to explore topological states of matter with great controllability. Specifically, we consider realizing the JC lattice in the context of superconducting quantum circuits, where each JC lattice site is constructed by a transmission line resonator (TLR) coupled to a two-level transmon qubit. We find that the dressed polariton states in the single-excitation manifold in each JC lattice site can simulate a spin-1/2 system. Particularly, synthetic SOC and Zeeman field for polaritons can be induced and manipulated by only engineering the coupling strength between neighboring resonators. Meanwhile, we show that, based on tunable synthetic SOC and Zeeman field, nodal-loop semimetal bands^{35–37} and topological polaritons can be realized and explored in the simulated JC lattice. Moreover, through calculating the topological winding number, we find that this tunable system has a rich topological phase diagram.

Our proposal to explore the topological states in the JC lattice system is different from previous ones based on the optical lattices.^{28–30} In particular, our proposal has a number of advantages. (i) Unlike ultracold atoms, polaritons are quasiparticles which are hybrids of photons and qubit excitations. Topological polaritons emerge from the topological structure of light–matter interaction, where photons and qubit excitations are topologically trivial by themselves, but combining together, they

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become hybrid topological states. Therefore, using polaritons for quantum simulation enriches our controlling methods—both photonic and atomic means take effects. (ii) The systematic parameters in JC lattice systems can be tuned at a single-site level, which allows us to generate a wide variety of SOC forms. (iii) The geometry of the JC lattice can be artificially designed, and lattice boundaries are easy to be created for observing topological edge states; thus, various topological lattice models and topological effects can be constructed and probed. (iv) The particle number putting in a JC lattice can be deterministically controlled. With such an advantage, we present a method using single-particle quantum dynamics to probe topological winding numbers and topological polariton edge states. (v) In the quantum optics platform, JC lattice systems previously have generated multiple important applications, including masers, lasers, photon transistors, and quantum information processors. Meanwhile, there are indeed several disadvantages in our proposed JC lattice system, such as the limited system size, parameter fluctuations, and decoherence. However, the essential physics of the simulated topological polariton states, such as the topological invariants and edge states, can still be detected under these realistic circumstances. Therefore, the topological JC lattice system in superconducting quantum circuits offers the possibility to develop functional topological spin quantum devices.

RESULTS

Janeys–Cummings lattice

The method for implementing a one-dimensional (1D) JC lattice in superconducting quantum circuits⁴ is as follows. As shown in Fig. 1a, every unit cell consists of a TLR resonantly coupled with a transmon, forming a JC model.³⁸ The neighboring TLRs are connected by a combination of a SQUID and a small inductor L in series, which can actually be regarded as the counterpart of a semitransparent mirror in the cavity QED system, allowing photons to hop across neighboring cavities (see the Methods section). As a result, setting $\hbar = 1$ hereafter, the system Hamiltonian of this JC lattice is

$$H_{JC} = \sum_{l=1}^N h_l + \sum_{l=1}^{N-1} J_l(t) (\hat{a}_l^\dagger \hat{a}_{l+1} + \text{h.c.}) + H_c, \quad (1)$$

where N is the number of the unit cells; $h_l = \omega_l (\sigma_l^+ \sigma_l^- + \hat{a}_l^\dagger \hat{a}_l) + g_l (\sigma_l^+ \hat{a}_l + \sigma_l^- \hat{a}_l^\dagger)$ is the JC-type interacting Hamiltonian in l th unit cell. The condition $g_l \ll \omega_l$ has to be met for justifying the JC coupling. $\sigma_l^+ = |e\rangle\langle g|$ and $\sigma_l^- = |g\rangle\langle e|$ are the raising and lowering operators of the l th transmon qubits. \hat{a}_l and \hat{a}_l^\dagger are the annihilation and creation operators of the photon in l th TLR. $J_l(t)$ is the inter-TLR hopping strengths between l th and $(l+1)$ th unit cells. Different from optical cavities, the time dependence of $J_l(t)$ here can be induced by adding a time-varying external magnetic flux threading through the SQUIDs (see Methods). $H_c = \sum_{l=1}^N g_l (\sigma_l^+ \hat{a}_l^\dagger + \sigma_l^- \hat{a}_l) + \sum_{l=1}^{N-1} J_l(t) (\hat{a}_l^\dagger \hat{a}_{l+1} + \text{h.c.})$ is the counter-rotating term that can be neglected by using the rotating wave approximation (RWA).

The lowest three eigenstates of the JC Hamiltonian h_l for the l th unit are $|0g\rangle_l$, $|\uparrow\rangle_l = (|0e\rangle_l + |1g\rangle_l)/\sqrt{2}$ and $|\downarrow\rangle_l = (|0e\rangle_l - |1g\rangle_l)/\sqrt{2}$, where $|ng\rangle_l$ and $|ne\rangle_l$ ($n = 0, 1, 2, \dots$) are the states containing n photons while the transmon is at the ground and excited state, respectively. Their eigenenergies are $E_{l,0g} = 0$ and $E_{l,1(\downarrow)} = \omega_l + (-)g_l$. Here, we exploit the two single-excitation eigenstates $|\uparrow\rangle_l$ and $|\downarrow\rangle_l$ to simulate the effective electronic spin-up and spin-down state. As each of the two states consists of a half “photon” and a half “atom”, they are regarded as a whole and termed as a “polariton”.

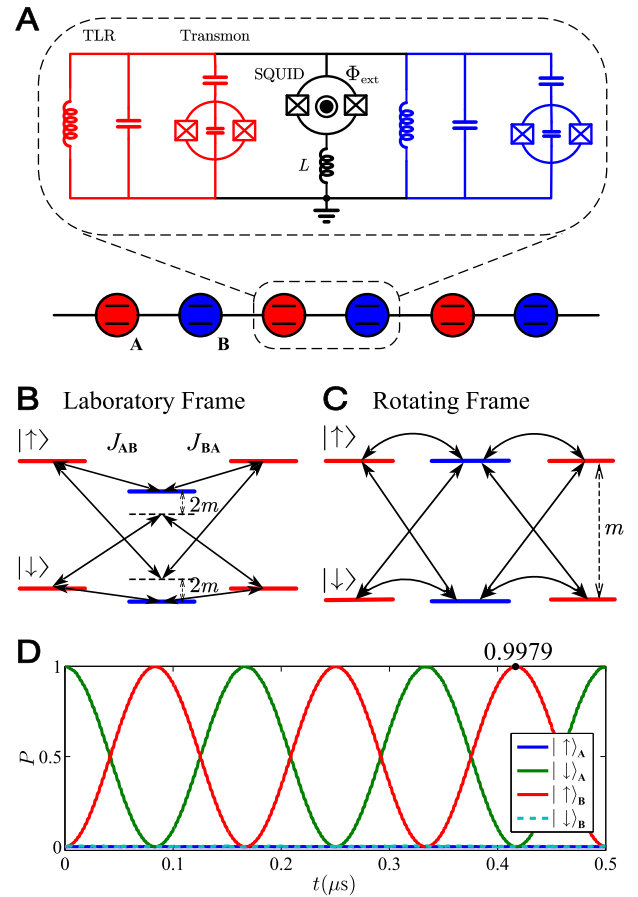


Fig. 1 The proposed superconducting circuit implementation of spin-1/2 lattice models. **a** The “spin-1/2” polariton lattice with two types of unit cells, A-type (red) and B-type (blue), arranged alternately. Each unit cell has two pseudo-spin-1/2 states simulated by the two single-excitation eigenstates of the JC model. The two types of unit cells are of the different qubit and photon eigenfrequencies and JC coupling strengths. The zoom-in figure details the equivalent superconducting circuits of two neighboring unit cells and their coupling circuit, which is a combination of a SQUID and an inductor L in series, to induce the tunable intercell photon hopping. **b** The resonant and detuning couplings of intercell spin states. Since the alternate red and blue unit cells arrangement, two sets of driving, J_{AB} and J_{BA} , have to be adopted to ensure the translation symmetry in the rotating frame defined by U . **c** The levels and designed hopping of the polariton lattice in the rotating frame, where the A-type and B-type unit cells can be treated as the same, so that the proposed circuit simulates a 1D spin-1/2 tight-binding lattice model. **d** The Rabi oscillation of a two-unit-cell system to justify the treatment of the proposed intercell coupling. The considered transition $|\uparrow\rangle_A \leftrightarrow |\downarrow\rangle_B$ is of the worst meeting the RWA requirement among the four possible transitions; thus, the fidelity obtained is the least one, but it still reaches a very high value of 0.9979 in the third Rabi cycle. All the numerical simulations are based on the Hamiltonian in Eq. 1 without RWA

Polaritonic spin–orbit coupling

We proceed to show that a spin-1/2 chain model with a tunable Zeeman field and SOC can be simulated with the JC lattice by only adjusting the pulse shape of the coupling strengths $J_l(t)$ between neighboring TLRs. First, since each cell contains two pseudo-spin states, there are totally four intercell neighboring hopping. In order to control each hopping separately, selective frequency addressing is employed (see Methods), i.e., we assign each of the four hopping with its unique hopping frequency. To achieve this, we adopt two sets of unit cells, A-type and B-type, which are

different in the sense that they have different eigenfrequencies and coupling strengths of the JC model. Then we arrange them in an alternate way, as shown in Fig. 1a. Setting started with an A-type one, when l is odd (even), $\omega_l = \omega_A(\omega_B)$ and $g_l = g_A(g_B)$. Then, based on the current experimental reaches,⁴ we set $\omega_A/2\pi = 6$ GHz, $\omega_B/2\pi = 5.65$ GHz, $g_A/2\pi = 300$ MHz, and $g_B/2\pi = 270$ MHz. With this, the energy intervals of the four hopping are $\{|E_{l,\alpha} - E_{l+1,\alpha'}|/2\pi\} = \{220, 320, 380, 920\}$ MHz with $\alpha, \alpha' \in \{\uparrow, \downarrow\}$. The differences between every two of them are no less than 20 times of the effective hopping strength $t_0/2\pi = 3$ MHz, thus they can be selectively addressed in frequency. Correspondingly, the $J_l(t)$ contain four tunes and can be written as

$$J_l(t) = \sum_{\alpha, \alpha'} 4t_{0,l,\alpha\alpha'} \cos(\omega_{l,\alpha\alpha'}^d t + s_{l,\alpha\alpha'} \varphi_{l,\alpha\alpha'}), \quad (2)$$

where

$$s_{l,\alpha\alpha'} = \text{sgn}(E_{l,\alpha} - E_{l+1,\alpha'}) \quad (3)$$

is the sign of the hopping phase, $4t_{0,l,\alpha\alpha'}$, $\omega_{l,\alpha\alpha'}^d$ and $s_{l,\alpha\alpha'}\varphi_{l,\alpha\alpha'}$ are the amplitudes, frequencies, and phases, corresponding to the hopping $|a\rangle_l \rightarrow |a'\rangle_{l+1}$, respectively. Note that due to the alternate arrangement, based on the definition above, the $J_l(t)$ s take only two different forms—when l is odd (even), $J_l = J_{AB}(J_{BA})$ where the subscript “AB” (“BA”) refers to that the A-type (B-type) unit cell is on the left. Experimentally, this time-dependent coupling strength $J_l(t)$ can be realized by adding external magnetic fluxes with both dc and ac components threading through the SQUIDs (see Methods). In this way, selective hopping can be induced, i.e., only when a frequency of the driving flux matches a particular hopping energy interval, that hopping can be triggered, otherwise it will not take into effect. Meanwhile, both the strengths and phases of the hopping can be controlled by the amplitudes and phases of the ac magnetic fluxes. However, only controlling these two is not enough for realizing the topological states; we still need to induce and adjust the spin splitting. Then we theoretically find out, and numerically prove that this spin splitting can be induced by just adding a detuning to the spin-flipped transition tunes. Concretely, while the spin-preserved transition frequencies are set as $\omega_{l,\alpha\alpha}^d = |E_{l,\alpha} - E_{l+1,\alpha}|$, we set the spin-flipped transition frequencies with a detuning $2m$ as $\omega_{l,\alpha\alpha'}^d = |E_{l,\alpha} - E_{l+1,\alpha'}| - 2m$, where $\alpha \neq \alpha'$, see Fig. 1b. Thus, in the rotating frame (explained later), there will be a spin splitting m for each cell, as shown in Fig. 1c.

We now show how the time-dependent coupling strength in Eq. 2 can induce a designable spin transition process in a certain rotating frame. First, we map the Hamiltonian in Eq. 1 into the single-excitation direct product subspace $\text{span}\{|0g, \dots, 0g, \alpha, 0g, \dots, 0g\rangle\}$. Hereafter, when there is no ambiguity, we use $|a\rangle_l$ to denote $|0g, \dots, 0g, \alpha, 0g, \dots, 0g\rangle$, and $|G\rangle$ to denote $|0g, \dots, 0g\rangle$. Then, we define a rotating frame by a unitary operator $U = \exp\{-i[\sum_l h_l - m(|\uparrow\rangle\langle\uparrow| - |\downarrow\rangle\langle\downarrow|)]t\}$, which leads the Hamiltonian in Eq. 1 to $H'_{JC} = U^\dagger H_{JC} U + i\dot{U}^\dagger U$. Neglecting the fast-rotating terms (see Methods), one obtains

$$H'_{JC} = \sum_l m \mathbf{S}_l^z + \sum_{l=1}^{N-1} \sum_{\alpha, \alpha'} \left(t_{0,l,\alpha\alpha'} e^{i\varphi_{l,\alpha\alpha'}} \hat{c}_{l,\alpha}^\dagger \hat{c}_{l+1,\alpha'} + \text{h.c.} \right), \quad (4)$$

where $\mathbf{S}_l^z = |\uparrow\rangle\langle\uparrow| - |\downarrow\rangle\langle\downarrow|$, $t_{0,l,\alpha\alpha'}$ is the effective coupling strength and $\hat{c}_{l,\alpha}^\dagger = |a\rangle_l \langle G|$ is the creation operator for a polariton with “spin” a in l th unit cell. As a result, the Hamiltonian (4) represents a general 1D spin-1/2 tight-binding lattice model with m , $t_{0,l,\alpha\alpha'}$, and $\varphi_{l,\alpha\alpha'}$ being the equivalent Zeeman energy, hopping strength, and hopping phase, respectively. These three variables can all be experimentally tuned in wide ranges by the frequencies, amplitudes, and phases of the external ac magnetic fluxes. Notably, the effective on-site potential m can be tuned as either positive or negative depending on the detuning direction.

Meanwhile, it is worth noticing that although we have introduced two kinds of unit cells (A-type and B-type) in the laboratory frame, by adjusting the coupling parameters J_{AB} and J_{BA} , the translation symmetry in the rotating frame is still preserved. In other words, the smallest repeating unit in the rotating frame contains only one unit cell as shown in Fig. 1c. By now, the adjustable spin-preserved tunneling ($\alpha = \alpha'$) and SOC terms ($\alpha \neq \alpha'$) can both be induced; hence, our superconducting quantum circuit setup can naturally be used to simulate a tunable SOC topological polariton insulator.

In order to justify the individual frequency addressing of the intercell transitions, we numerically simulate the dynamics of a system containing only two unit cells, with an A-type one on the left and a B-type one on the right. We test every hopping of the four transitions $\{|a\rangle_A \leftrightarrow |a'\rangle_B\}$ one by one, adding only one corresponding frequency $\omega_{1,\alpha\alpha'}^d$ in $J_{AB}(t)$. As expected, when we pick out a resonant frequency $\omega_{1,\alpha\alpha'}^d$ with $m = 0$, there will be a Rabi oscillation between the two corresponding target states. One example of these Rabi oscillations was shown in Fig. 1d, which is of the least fidelity among the four. Even in this worst case, and in the third Rabi cycle, the fidelity still reaches a very high value of 0.9979, which justifies the RWA. In addition, our numerical simulation also shows that, in the presence of the unmatched driving, all the initial nontarget states remain almost unchanged, thus justifying our individual frequency-addressing method.

Nodal-loop topological polaritons

Our protocol provides a tunable platform using polaritons to study topological matters. Here, we take the nodal-loop semimetal as an application sample to demonstrate how to simulate a specific condensed matter model in our proposed setup. To fit our simulation setup, we reform the Hamiltonian of the original 3D nodal-loop model in ref.³⁷, without losing any topological properties. First, we relabel the coordinates to set the hopping terms with SOC to be along the x axis. Second, we consider the Fourier transformations along y and z directions with quasi-momenta k_y and k_z and treat them as system parameters. Then, according to Eq. 4, we can simulate the 3D nodal-loop model in our 1D system with the other two dimensions being the parametric dimensions. In this direct simulation, to engineer the four transitions between different polariton states, we need four different tunes in the intercell coupling strength $J_l^{\text{nod}}(t)$.

The above implementation can further be simplified as the following. We first make a unitary transformation to the original Hamiltonian, so that $H_{\text{nod}} = V^\dagger H_{\text{ori}} V$, where $V = \sum_{l=1}^N (-i)^{l-1} I_l$ with $I_l = |\uparrow\rangle\langle\uparrow| + |\downarrow\rangle\langle\downarrow|$. The transformed 1D lattice Hamiltonian from Eq. 4, without losing any physical properties, is

$$H_{\text{nod}} = \sum_{l=1}^N m'(k_y, k_z) \mathbf{S}_l^z + \sum_{l=1}^{N-1} \sum_{\alpha, \alpha'} i t_0' \hat{c}_{l,\alpha}^\dagger \hat{c}_{l+1,\uparrow} - i t_0' \hat{c}_{l,\downarrow}^\dagger \hat{c}_{l+1,\downarrow} + i t_0' \hat{c}_{l,\uparrow}^\dagger \hat{c}_{l+1,\downarrow} - i t_0' \hat{c}_{l,\downarrow}^\dagger \hat{c}_{l+1,\uparrow} + \text{h.c.}, \quad (5)$$

where $m'(k_y, k_z) = M + 2d(\cos k_y + \cos k_z)$ with M being the effective Zeeman energy and d being the effective hopping energy along y and z directions. Then, we set the parameters of the JC model to be $\omega_A = \omega_B = 2\pi \times 6$ GHz, $g_A/2\pi = 200$ MHz, and $g_B/2\pi = 100$ MHz. Correspondingly, the coupling strengths can be set to contain only two tunes as

$$J_l^{\text{nod}}(t) = 4t_0 \cos \left[\omega_1^d t + \frac{(-1)^{l+1}\pi}{2} \right] + 4t_0 \cos \left(\omega_2^d t + \frac{\pi}{2} \right), \quad (6)$$

where $\omega_1^d/2\pi = 100$ MHz and $\omega_2^d/2\pi = (300 - 2m)$ MHz. In this way, when transforming into the rotating frame of U and applying the RWA, the Hamiltonian (1) of our JC-lattice system will take the

form of Eq. 5, which accomplishes the quantum simulation of the topological nodal-loop semimetals (see Methods).

Experimentally, one can choose $t_0/2\pi = 3$ MHz and set the detuning within the range of $m \in 2\pi \times [-20, 20]$ MHz, such that any value of the variable tunes $\omega_{\alpha\alpha'}^d$ ($\alpha \neq \alpha'$) still maintains a frequency difference no less than $20t_0$ away from other tunes, and thus the crosstalk caused by unwanted tunes will be negligible. Last, for testing the validity of our theoretical protocol, numerical simulations will be given in the Experimental detection section after a brief introduction of the characteristics of the topological nodal-loop polaritons.

Characteristics of the topological polaritons

To investigate the bulk characteristic of the topological nodal-loop polaritons, we first consider the periodic boundary condition to obtain the Hamiltonian in Eq. 5 in the momentum space as

$$H = b_y \mathcal{S}^y + b_z \mathcal{S}^z, \quad (7)$$

where $b_y = -2t'_0 \cos k_x$, $b_z = 2t'_0 \sin k_x + m'(k_y, k_z)$ and $\mathcal{S}^y = i|\downarrow\rangle\langle\uparrow| - i|\uparrow\rangle\langle\downarrow|$. This system has two energy bands

$$E = \pm \sqrt{b_y^2 + b_z^2}, \quad (8)$$

which will touch when $E=0$. The touching points form closed lines, the so-called nodal loops, in momentum space as shown in Fig. 2a. These two loops appear in the $k_x = \pi/2$ and/or $k_x = -\pi/2$ planes. By fixing $k_x = \pm\pi/2$, we plot these two energy bands in the k_y - k_z space as shown in Fig. 2b, c, where one can see that the touching is right along the nodal loops.

The topological index characterizing each nodal loop is a winding number defined as

$$\nu(k_y, k_z) = \frac{1}{2\pi} \int_{-\pi}^{\pi} dk_x \mathbf{v} \times \partial_{k_x} \mathbf{v} \\ = \frac{1}{2} [\text{sgn}(m' + 2t'_0) - \text{sgn}(m' - 2t'_0)], \quad (9)$$

where $\mathbf{v} = (v_y, v_z) = (b_y, b_z)/\sqrt{b_y^2 + b_z^2}$. This shows that the quantized winding number is either 1 or 0, corresponding to the cases whether a nodal loop is enclosing the straight line, which is along the k_x direction of the fixed (k_x, k_y) point, or not.^{37,38} Hence, the nodal loops divide the k_y - k_z space into regions with different winding numbers, as shown in Fig. 2d. For all straight lines along k_x inside the nodal loop, each of them can be regarded as being corresponding to a topological 1D-gapped subsystem with winding number 1.

Two striking topological characters of the nodal-loop semimetal are the zero-energy edge states inside the energy gap and their corresponding edge states. We take a slice of $k_y = 0$, indicated by the green line in Fig. 2d, as an example to plot the energy spectrum with various k_z for a finite chain with $N = 20$, as shown in Fig. 2e. The numerical result shows that there are two mid-gap degenerated zero-energy modes appearing in the range of $\nu = 1$ (the red area in Fig. 2d). The quantum states corresponding to the two mid-gap energies are edge states localized in the left and right end of the lattice, respectively. When N is large enough for ignoring the finite-size effects, these wave functions can be

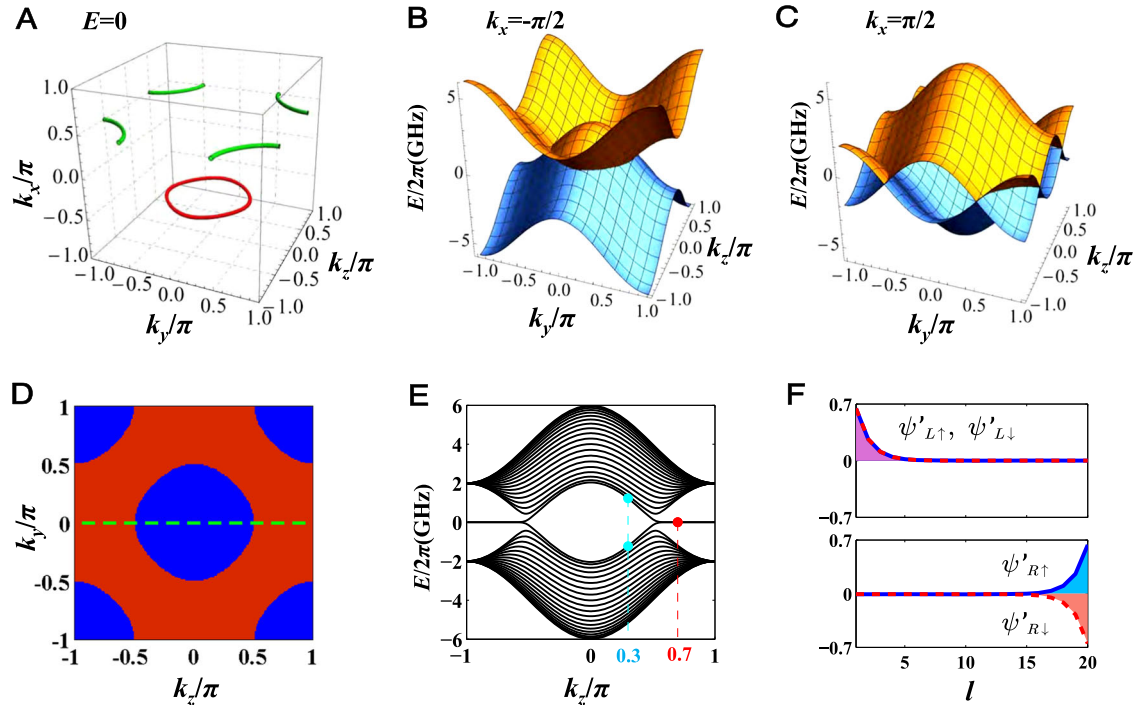


Fig. 2 The band structure and topological characteristics of the simulated nodal-loop semimetal. **a** The two loops (lines of $E = 0$) where the two bands touch in the momentum space. There is one in the $k_x = -\pi/2$ plane (colored red) and one in the $k_x = \pi/2$ plane (colored green). The parameters for plotting are $M/t_0 = 0$ and $d/t_0 = 1$. Energy bands under the confines of **b** $k_x = -\pi/2$ and **c** $k_x = \pi/2$. They touch each other along nodal loops. **d** The winding number altering over k_y and k_z . The red and blue regions are of winding number $\nu = 1$ and $\nu = 0$, respectively. **e** Numerical calculation of the energy bands of a 1D 20-unit-cell lattice along the x direction in open boundary condition. The energies are altering over k_z with confining $k_y = 0$ (the dashed green line in **d**). The two light blue dots at $k_z = 0.3\pi$ refer to two topological trivial states used for comparison in later discussion. The two red dots (overlapped) at $k_z = 0.7\pi$ mark two in-gap zero-energy levels accompanied with two edge states whose wave functions are plotted in **f**, where the blue and dashed red lines plot the probability $|\uparrow\rangle$ and $|\downarrow\rangle$ components of the numerically calculated wave functions divided by a phase factor i^{l-1} , i.e., $\psi'_{L(R)\uparrow(\downarrow)}(l) = \psi_{L(R)\uparrow(\downarrow)}^{\text{num}}(l)/i^{l-1}$

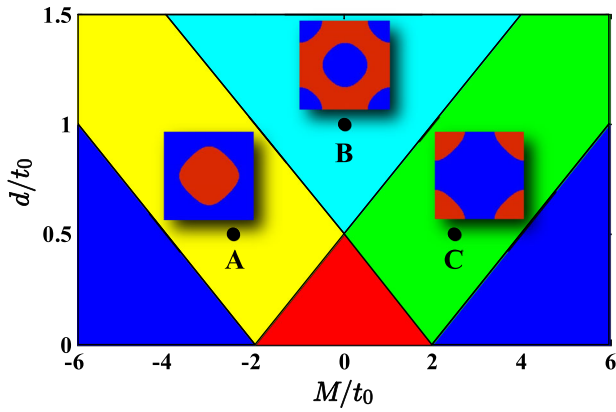


Fig. 3 Phase diagram of nodal-loop semimetal bands with the corresponding winding number configurations. Each color denotes a different phase. The dark blue region is of a trivial gapped phase without nodal loops in momentum space; the yellow and green region is of only one nodal loop in the k_y - k_z plane; the light blue region is of two nodal loops in the k_y - k_z plane; the red region is of a nontrivial gapped phase without nodal loops but with a winding number $\nu=1$ in the whole k_y - k_z plane. The three insets are the configuration of the winding number in the k_y - k_z plane (the red regions are of $\nu=1$, while blue regions are of $\nu=0$), corresponding to points A $(-2.5, 0.5)$, B $(0, 1)$, and C $(2.5, 0.5)$, respectively

expressed as

$$\psi_L = \sum_{l=1}^N i^{l-1} A e^{-\lambda(l-1)} (|\uparrow\rangle_l + |\downarrow\rangle_l), \quad (10a)$$

$$\psi_R = \sum_{l=1}^N i^{l-1} A e^{-\lambda(N-l)} (|\uparrow\rangle_l - |\downarrow\rangle_l), \quad (10b)$$

where $A = \sqrt{(1-q^2)/2(1-q^{2N})}$ and $\lambda = \ln(1/q)$ with $q = \tan(m\pi/8t_0)$. The phase factor i^{l-1} stems from the unitary transformation V to get the nodal-loop Hamiltonian in Eq. 5 that simplified our simulation, from the original Hamiltonian in Eq. 4. These analytical results of wave functions are in very good agreement with the numerical results for $N=20$, as shown in Fig. 2f, thus justifying that we can use the JC lattice of an experimentally capable size to simulate the topological features.

There are several phases in our simulated Hamiltonian, where the phase transition is indicated by the emerging or vanishing of the nodal loops. Inferring from Eq. 8, the critical conditions are obtained as $k_x = -\pi/2$ and $k_x = \pi/2$, which are corresponding to one nodal loop in each of the two regions of $-2t'_0 - 4d < M < -2t'_0 + 4d$ and $2t'_0 - 4d < M < 2t'_0 + 4d$ with $d > 0$ in the M - d plane. Therefore, in the area where the two regions overlap, there are two nodal loops. But in the area outside these two regions, there is no nodal loop, so that the whole Brillouin zone will be in a purely trivial or nontrivial phase. Consequently, there are totally five different phases of different winding number configurations $\nu(k_y, k_z)$ in the M - d plane, as shown in Fig. 3. According to the chosen parameters, t_0 and m , the area $\{M, d | -6t'_0 + 4d < M < 6t'_0 - 4d\}$ will include all the five phases in the phase diagram of Fig. 3, i.e., all the possible phases can be simulated in our proposed system.

Experimental detection methods

Polaritonic topological edge state detection. According to Eqs. 10a and 10b, or as shown in Fig. 2f, the polariton in the left or right edge state is maximally distributed in the leftmost and rightmost JC lattice sites. Their internal spins are in the superposition states $(|\uparrow\rangle_l + |\downarrow\rangle_l)/\sqrt{2}$ and $(|\uparrow\rangle_l - |\downarrow\rangle_l)/\sqrt{2}$, respectively. One can find that, in the beginning, the left and right polaritonic edge states only have qubit excitation and photon components, respectively.

Taking the detection of the left edge state as an example, initially, the polariton in the leftmost JC lattice site is prepared into $|0e\rangle_1$, i.e., the leftmost qubit (resonator) has been prepared in the excited (vacuum) state. The qubits and resonators in the other sites are prepared into the ground and vacuum states, which means the initial systematic state is $|\psi(t=0)\rangle = |0e\rangle_1 |0g\rangle_2 \cdots |0g\rangle_N$. After that, we let the above initial state evolve for a time of about $0.5 \mu\text{s}$. If the JC lattice is in the topological nontrivial phase supporting the left edge states, the final density distribution of the polaritons will maximally populate the leftmost site. The reason is that the initial state $|\psi(t=0)\rangle$ has a large overlap with the left edge state. It will evolve mainly via the edge state wave packet and is maximally localized in the leftmost site. While if the system is in the topological trivial phase and has no edge states, the initial state will be a superposition of different bulk states. The final density distribution will not have maximal distribution in the leftmost site. Similarly, one can also prepare the JC lattice into $|\psi(t=0)\rangle = |0g\rangle_1 \cdots |0g\rangle_{N-1} |1g\rangle_N$ and detect³⁹ the right polaritonic topological edge state based on observing its time evolution.

In Fig. 4a, b, we have numerically calculated the time evolution of the polaritonic density when the JC lattice is in the topological trivial and nontrivial nodal-loop semimetal phases, respectively. For the trivial case, the wave packet has a ballistic spread versus time, which is a typical feature of bulk Bloch state. It shows that there is no edge state localization and the system is in a topological trivial phase. For the nontrivial case, the density of the polaritons will always maximally localize in the leftmost JC lattice site, which indicates the existence of the left topological edge states, demonstrating that the system is in a topological nontrivial phase. The time evolution of the qubit excitation and the photon population for the topological nontrivial case is also numerically calculated in Fig. 4c, d, which shows that the localized qubit excitation and the photon in the leftmost site have a Rabi-like oscillation feature inherited from the JC model.

Polaritonic topological invariant detection. Another important hallmark for a topological nontrivial nodal-loop polaritonic semimetal phase is the nontrivial polaritonic topological winding number. Here we show that such a polaritonic topological invariant can also be dynamically detected. Our method is based on a previous work which shows that the topological winding number rotted in the momentum space can be detected through measuring the dynamical chiral center in the real space.⁴⁰ The chiral operator for our topological polaritonic model is $\mathbf{S}_l^x = |\uparrow\rangle\langle\downarrow| + |\downarrow\rangle\langle\uparrow|$. Then the chiral center operator for the JC lattice is defined as $\hat{P}_d = \sum_{l=1}^N l \mathbf{S}_l^x$. The polaritonic topological winding number can be related with the time-averaged dynamical chiral center associated with the single-polariton dynamics, i.e.,

$$\nu = \lim_{T \rightarrow \infty} \frac{2}{T} \int_0^T dt \langle \psi_c(t) | \hat{P}_d | \psi_c(t) \rangle, \quad (11)$$

where T is the evolution time, $|\psi_c(t)\rangle = \exp(-iH_{\text{nod}}t)|\psi_c(0)\rangle$ is the time evolution of the initial single-polariton state $|\psi_c(0)\rangle = |0g\rangle_1 \cdots |\uparrow\rangle_{[N/2]} \cdots |0g\rangle_N$, where one of the middle JC lattice site has been put into one polariton, with its spin prepared in the state $|\uparrow\rangle$.

In Fig. 5a, b, we have numerically calculated the dynamical chiral center $\bar{P}_d(t) = \langle \psi_c(t) | \hat{P}_d | \psi_c(t) \rangle$ for topological trivial and nontrivial cases, respectively. According to Eq. 11, one can find that the topological winding number is equal to twice the oscillation center of $\bar{P}_d(t)$. As shown in Fig. 5a, $\bar{P}_d(t)$ oscillates around the average value 0, which gives the polaritonic topological winding number $\nu=0$. In contrast, the result for the topological nontrivial case in Fig. 5b shows that $\bar{P}_d(t)$ oscillates around 0.5, which yields the polaritonic topological winding number $\nu=1$. Experimentally, the states of qubits and resonators can be measured with a fidelity higher than 0.99 in

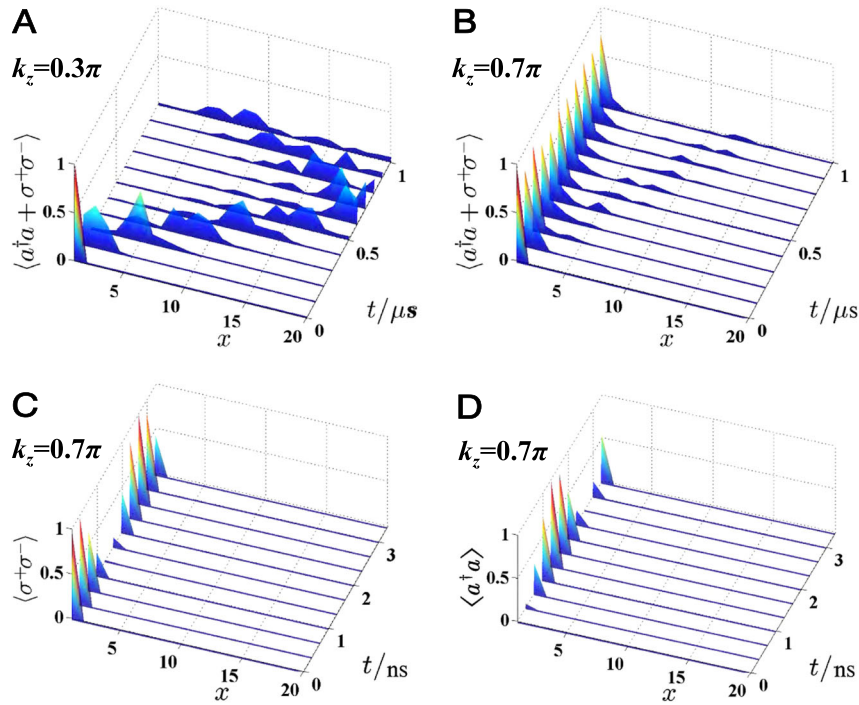


Fig. 4 Dynamical detection of polaritonic topological edge states. Time evolution of a polaritonic density distribution $\langle \sigma^+ \sigma^- + \hat{a}^\dagger \hat{a} \rangle$ when the JC lattice is in **a** topological trivial phase of $k_z = 0.3\pi$ and **b** topological nontrivial phase of $k_z = 0.7\pi$ (see Fig. 2e). Time evolution of **c** qubit excitation distribution $\langle \sigma_1^+ \sigma_1^- \rangle$ and **d** photon distribution $\langle \hat{a}_1^\dagger \hat{a}_1 \rangle$ for the topological nontrivial case. All the numerical simulations are based on the Hamiltonian in Eq. 1 without RWA. Other parameters are the same as those in Fig. 2e

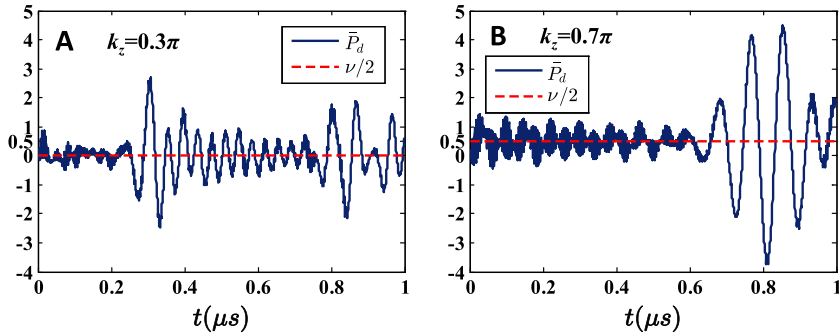


Fig. 5 Dynamical detection of polaritonic topological invariants. Time evolution of the chiral center \bar{P}_d when the JC lattice is in **a** the topological trivial phase of $k_z = 0.3\pi$ and **b** nontrivial phase of $k_z = 0.7\pi$ (see Fig. 2e). The red dashed line denotes the oscillation center. All the numerical simulations are based on the Hamiltonian in Eq. 1 without RWA. Other parameters are chosen the same as those in Fig. 4

superconducting circuits. In our case, one only needs to measure the qubit excitation and photon populations for getting their imbalance and deriving the chiral center, without requiring full quantum state tomography. In this way, the topological winding number can be easily and unambiguously detected based on monitoring single-polariton quantum dynamics in a JC lattice.

DISCUSSION

We further investigate how our construction and detection methods are influenced by the quantum decoherence effects. The Lindblad master equation is adopted to take three main decoherence factors, including the losses of the photon and the decay and dephasing of the transmon into account. The Lindblad master equation of our system can be written as

$$\dot{\rho} = -i[H_{JC}, \rho] + \sum_{l=1}^N \sum_{i=1}^3 \gamma \left(\Gamma_{l,i} \rho \Gamma_{l,i}^\dagger - \frac{1}{2} \{ \Gamma_{l,i}^\dagger \Gamma_{l,i}, \rho \} \right), \quad (12)$$

where ρ is the density operator of the whole system, γ is the decay rate or noise strength which are set to be the same here, $\Gamma_{l,1} = a_l$, $\Gamma_{l,2} = \sigma_l^-$ and $\Gamma_{l,3} = \sigma_l^z$ are the photon-loss, transmon-loss, and the transmon-dephasing operators in the l th lattice, respectively. In Fig. 6a, we plot the edge-site population $P_1(t) = \text{trace}[\rho(t)(a_1^\dagger a_1 + \sigma_1^+ \sigma_1^-)]$ after $0.5 \mu\text{s}$ and the oscillation center $\nu/2$ of the trivial and nontrivial cases for different decay rates. It shows that the edge state population and the chiral center smoothly decrease when the decay rate increases. However, our detection method can tolerate the decay rate up to the order of $2\pi \times 100 \text{ kHz}$, while the typical decay rates are only $2\pi \times 5 \text{ kHz}$.

We also investigate the minimum sites that are needed for the experimental detection of the oscillation center. In the presence of the decay rates of $2\pi \times 5 \text{ kHz}$, we plot the oscillation center for a lattice with a different number of sites in both the trivial and nontrivial cases, as shown in Fig. 6b. It shows that the four-site case corresponding to an oscillation center value of about 0.40, is

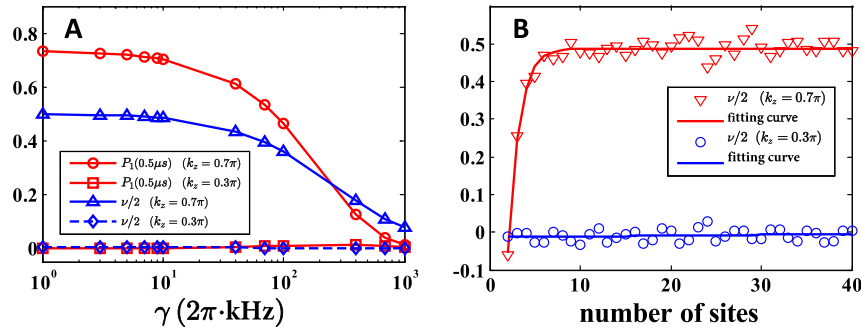


Fig. 6 The influences of the decoherence and the lattice site on the detection of the topological effects. **a** The edge-site population $P_1(t) = \text{trace}[\rho(t)(a_1^\dagger a_1 + \sigma_1^+ \sigma_1^-)]$ at $0.5 \mu\text{s}$ and the oscillation center $\nu/2$ of the topological trivial ($k_z = 0.3\pi$) and nontrivial ($k_z = 0.7\pi$) cases for different decay rate γ . **b** The oscillation center of the topological trivial ($k_z = 0.3\pi$) and nontrivial ($k_z = 0.7\pi$) cases varying over the number of lattice sites. The decay rates are all set to be $2\pi \times 5 \text{ kHz}$. All the numerical simulations are based on the Hamiltonian in Eq. 1 without RWA. Other parameters are chosen the same as those in Fig. 4

big enough for distinguishing the topologically trivial and nontrivial cases with current state-of-the-art technologies.

To conclude, we have introduced the concept of topological states into the JC lattice, which is one of the most important building blocks in quantum optics and quantum information processing. We have studied the topological structure of light-matter interaction and shown that SOC physics and topological polaritons can be pursued in the JC lattice. Different from synthetic topological states in ultracold atomic, photonic, and acoustic systems, topological polaritons are topological superposition states of photons and qubits. Tunable synthetic polaritonic SOC is induced by engineering the JC lattice couplings, which provides the basic ingredient for realizing spinful topological states of matter. We have also provided a method using single-particle quantum dynamics in real space to directly observe the polaritonic topological edge states and topological invariants.

Our work has a broad generalization and opens the door for exploring spinful topological states of matter using polaritons in the JC lattice system. (i) In addition to mimicking spin-1/2, polariton states in multiple-excitation manifold have an extended spin degree of freedom and can also be used to mimic a larger spin. It is challenging to realize SOC for the larger spin case in solid-state materials and ultracold atoms. However, the method proposed in our work can be generalized to realize synthetic SOC for large-spin polaritons. This allows us to explore a large variety of topological states, including triple-point topological states of matter,⁴¹ (ii) polariton states, the eigenstates of a JC model, can be referred as a synthetic dimension, where each pair of states mimic a spin in a spin-lattice model and the coupling between polariton states provides the hopping in the synthetic dimension. With such a synthetic dimension, high-dimensional topological states of matter can be explored in a low-dimensional JC lattice, including topological states beyond three dimensions,^{42,43} (iii) polaritons have a tunable strong nonlinear interaction, which allows us to study polaritonic fractional topological states of matter;⁴⁴ (iv) besides a superconducting circuit, JC lattices can also be realized in many quantum optical systems, including trapped ions,² cavity quantum electrodynamics,³ nanoscopic lattices,⁶ optomechanical systems,⁷ and so on.

METHODS

SQUID induced time-dependent photon hopping

We here present how to induce the time-dependent photon-hopping strength between two TLRs. As shown in Fig. 1a, neighboring TLRs are connected by a common SQUID and an inductor L in series and then to the ground, which actually serves as a single Josephson junction (JJ) but with an effective Josephson inductance tunable by the external flux. Concretely, by applying an external magnetic flux $\Phi_{\text{ext}} = \Phi_{\text{dc}} + \Phi_{\text{ac}}$ threading through

the SQUID, when $\Phi_{\text{ac}} \ll \phi_0$ with ϕ_0 being the reduced flux quanta, the effective inductance of the SQUID reads⁴⁵

$$L_S(\Phi_{\text{ext}}) = \frac{\phi_0}{2l_c \cos(\Phi_{\text{ext}}/2\phi_0)}, \quad (13)$$

where l_c is the shared critical current of the two JJ in each SQUID. On the other hand, comparing with the inductance of the TLRs, both the SQUID and the inductor L have relatively small inductances; thus, there is a voltage node but a current peak at both ends of each TLR. For these boundary conditions, after the conventional quantization of the TLRs,^{36,45} the flux density and the charge density wave function of the lowest-energy mode in l th TLR can be expressed as

$$\hat{\phi}_l(x_l, t) = \frac{\sqrt{\omega_l L_l}}{d_l} \cos\left(\frac{\pi}{d_l} x_l\right) [\hat{a}_l^\dagger(t) + \hat{a}_l(t)], \quad (14a)$$

$$\hat{q}_l(x_l, t) = i \frac{\sqrt{\omega_l L_l}}{d_l} \sin\left(\frac{\pi}{d_l} x_l\right) [\hat{a}_l^\dagger(t) - \hat{a}_l(t)], \quad (14b)$$

where $\omega_l = \pi/\sqrt{L_l C_l}$ is the frequency of the photon with L_l and C_l being the inductance and capacitance, respectively. d_l is the length and x_l is the coordinate of the l th TLR. Meanwhile, because of the relatively low impedances of the SQUID and the inductor, the currents from the ends of every TLR will flow directly through them to the ground, without crossing to their neighboring TLRs. Hence, the interaction Hamiltonian between the l th and $(l+1)$ th TLR is just the summation energy of the SQUID and the inductor L

$$\begin{aligned} H_{\text{int}}^l &= \frac{1}{2}(L_S + L)(I_l^i + I_{l+1}^e)^2 \\ &= \sum_{j=1}^{l+1} \frac{\omega_j}{2L_j} (L_S + L) (\hat{a}_j^\dagger + \hat{a}_j)^2 \\ &\quad - \sqrt{\frac{\omega_l \omega_{l+1}}{L_l L_{l+1}}} (L_S + L) (\hat{a}_l^\dagger + \hat{a}_l) (\hat{a}_{l+1}^\dagger + \hat{a}_{l+1}), \end{aligned} \quad (15)$$

where $I_l^{i(e)} = \hat{\phi}_l(x_l, t) d_l / L_l |_{x_l=d_l(0)}$ is the current of the right (left)-end of the l th TLR. Moreover, if we set

$$\Phi_{\text{ac}} = 2\phi_0 \arccos \frac{-1}{1 + \sum_j^{n'} \Omega_j [\cos(\omega_j^d t + \varphi_j) + 1]}, \quad (16a)$$

$$t_{0j} = \frac{\phi_0 \Omega_j}{8l_c} \sqrt{\frac{\omega_l \omega_{l+1}}{L_l L_{l+1}}}, \quad (16b)$$

$$L = \frac{\phi_0 \sum_j^{n'} (\Omega_j + 1)}{2l_c}, \quad (16c)$$

where n' is the number of the tunes in Φ_{ac} , $\Phi_{\text{dc}} = 4\pi n'' \phi_0$ where n'' is an arbitrary positive integer, and choose resonant or detuned frequencies of ω_j after the RWA, we obtain

$$H_{\text{int}}^l = 4 \sum_j^{n'} t_{0j} \cos(\omega_j^d t + \varphi_j) (\hat{a}_l^\dagger \hat{a}_{l+1} + \text{h.c.}), \quad (17)$$

which correctly meets the form of Eqs. 1 and 2. This equation can be interpreted as describing the photon hopping between neighboring unit

cells, which means that the SQUID-L combination can actually serve as a counterpart of the semitransparent mirror in the optical cavity system.

Frequency-addressing control

We now show how the selective control of individual hopping in the JC lattice can be achieved by adjusting $J_l(t)$ in Eq. 1 via the ac flux. We first map the Hamiltonian in Eq. 1 into the single-excitation subspace span $\{|a\rangle_l\}_{l=1,2,\dots,N; a=\uparrow,\downarrow}$ and get

$$H_{JC} = \sum_{l,a} E_{l,a} \hat{a}_{l,a}^\dagger \hat{c}_{l,a} + \frac{1}{2} \sum_{l=1}^{N-1} \sum_{a,a'} J_l(t) \hat{c}_{l,a}^\dagger \hat{c}_{l+1,a'} + \text{h.c.}, \quad (18)$$

where $\hat{c}_{l,a}^\dagger = |a\rangle_l \langle G|$. Then in every $J_l(t)$, we add four tunes, corresponding to the four intercell hopping, cf. Eq. 17, each of which contains its independent tunable amplitude, frequency, and phase as

$$J_l(t) = \sum_{a,a'} 4t_{0,laa'} \cos(\omega_{laa'}^d t + s_{laa'} \varphi_{laa'}), \quad (19)$$

where $s_{laa'}$ is defined in Eq. 3 being the sign of each phase and the frequencies are set as

$$\omega_{laa'}^d = \begin{cases} |E_{l,a} - E_{l+1,a'}| & (a = a'), \\ |E_{l,a} - E_{l+1,a'}| - 2m & (a \neq a'), \end{cases} \quad (20)$$

where $2m$ is a detuning. By now, the form of $J_l(t)$ is determined, leaving m , $t_{0,laa'}$, and $\varphi_{laa'}$ to be chosen arbitrarily depending on the topological insulator model one simulates. Eventually, the target topological insulator model is in the rotating frame transformed by

$$U = e^{-i[\sum_l h_l - m(\uparrow\uparrow) + (1-\downarrow\downarrow)]t}, \quad (21)$$

After the picture transformation $H'_{JC} = U^\dagger H_{JC} U + iU^\dagger \dot{U}$, one gets

$$H'_{JC} = \sum_{l=1}^N m \mathbf{S}_l^\dagger + \sum_{l=1}^{N-1} \sum_{a,a'} t_{0,laa'} [e^{i(\omega_{laa'} t + s_{laa'} \varphi_{laa'})} + e^{-i(\omega_{laa'} t + s_{laa'} \varphi_{laa'})}] \times \begin{bmatrix} e^{i(E_{l+1,1} - E_{l,1})t} \hat{c}_{l,1}^\dagger \hat{c}_{l+1,1}^\dagger + e^{i(E_{l+1,1} - E_{l,1})t} \hat{c}_{l,1}^\dagger \hat{c}_{l+1,\downarrow}^\dagger \\ + e^{i(E_{l+1,1} - E_{l,1} - 2m)t} \hat{c}_{l,1}^\dagger \hat{c}_{l+1,\uparrow}^\dagger + e^{i(E_{l+1,1} - E_{l,1} + 2m)t} \hat{c}_{l,1}^\dagger \hat{c}_{l+1,\downarrow}^\dagger \end{bmatrix} + \text{h.c.}, \quad (22)$$

where $\mathbf{S}_l^\dagger = |\uparrow\rangle\langle\uparrow| - |\downarrow\rangle\langle\downarrow|$ and $\hat{c}_{l,a}^\dagger = |a\rangle_l \langle G|$. After doing the multiplication in this equation, the four resonant terms and their hermitian conjugates will be absent of time t (frequency addressing). Meanwhile, if the conditions $\{t_{0,laa'}\}_{a,a'=\uparrow,\downarrow} \ll \{\omega_{laa'}, \omega_{laa'} - \omega_{l\beta\beta'}\}_{a,a',\beta,\beta'=\uparrow,\downarrow; \omega_{laa'} \neq \omega_{l\beta\beta'}}$ are satisfied, all the other terms are a fast-rotating term that can be dropped within RWA. As a result, we obtain the tight-binding model with a tunable SOC in Eq. 4.

Simplified method of implementation

We here explain why the four hopping terms in the nodal-loop semimetal model can be induced by the coupling strength in Eq. 6 using only two tunes. First, when we set parameters of the JC model to be $\omega_A = \omega_B = 2\pi \times 6$ GHz, $g_A/2\pi = 200$ MHz, and $g_B/2\pi = 100$ MHz, the energy intervals of the four hopping overlap and reduce into two sets of intervals, i.e., a spin-conserved hopping interval

$$|E_{l,\uparrow} - E_{l+1,\uparrow}| = |E_{l,\downarrow} - E_{l+1,\downarrow}| = 2\pi \times 100 \text{ MHz}, \quad (23)$$

and a spin-flipped hopping interval

$$|E_{l,\uparrow} - E_{l+1,\downarrow}| = |E_{l,\downarrow} - E_{l+1,\uparrow}| = 2\pi \times 300 \text{ MHz}. \quad (24)$$

Therefore, in this way, one tune in $J_l^{\text{nod}}(t)$ can induce two hopping, while the two spin-conserved hopping still remain being controlled separated from the two spin-flipped hopping, thus they can be of different detuning.

Second, note that within this setting, according to the definition in Eq. 3, we have $s_{l\uparrow\uparrow} = -s_{l\downarrow\downarrow}$ and $s_{l\uparrow\downarrow} = -s_{l\downarrow\uparrow}$; thus, both the spin-conserved hopping terms and spin-flipped hopping terms can be induced by the same tune, but with opposite signs, i.e., i and $-i$. This is possible because the RWA selects different terms. These opposite signs are ideal for realizing the wanted SOC in our protocol. Anyway, this is only possible when we do the unitary transformation V to transform the hopping phase from the original 1 and -1 into the pure imaginary numbers i and $-i$. For the former case, one still has to use two tunes and induce them separately. Therefore, this transformation simplifies our simulation.

DATA AVAILABILITY

Data and analysis code used in this study are available from the corresponding authors on request.

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AUTHOR CONTRIBUTIONS

Z.Y.X. proposed the project. Z.Y.X., F.M., D.W.Z., and F.L.G. developed and designed the theoretical framework. F.L.G. performed the calculations with input from all the other authors. All authors discussed the contents and wrote the paper.

ADDITIONAL INFORMATION

Competing interests: The authors declare that they have no competing interests.

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