Hopf insulators and their topologically protected surface states

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Three-dimensional (3D) topological insulators in general need to be protected by certain kinds of symmetries other than the presumed $U(1)$ charge conservation. A peculiar exception is the Hopf insulators which are 3D topological insulators characterized by an integer Hopf index. To demonstrate the existence and physical relevance of the Hopf insulators, we construct a class of tight-binding model Hamiltonians which realize all kinds of Hopf insulators with arbitrary integer Hopf index. These Hopf insulator phases have topologically protected surface states and we numerically demonstrate the robustness of these topologically protected states under general random perturbations without any symmetry other than the $U(1)$ charge conservation that is implicit in all kinds of topological insulators.

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Topological phases of matter may be divided into two classes: the intrinsic ones and the symmetry protected ones. Symmetry protected topological (SPT) phases are gapped quantum phases that are protected by symmetries of the Hamiltonian and cannot be smoothly connected to the trivial phases under perturbations that respect the same kind of symmetries. Intrinsic topological (IT) phases, on the other hand, do not require symmetry protection and are topologically stable under arbitrary perturbations. Unlike SPT phases, IT phases may have exotic excitations bearing fractional or even non-Abelian statistics in the bulk. Fractional quantum Hall states and spin liquids belong to these IT phases. Remarkable examples of the SPT phases include the well known 2D and 3D topological insulators and superconductors protected by time reversal symmetry, and the Hal-dane phase of the spin-1 chain protected by the $SO(3)$ spin rotational symmetry. For interacting bosonic systems with on-site symmetry $G$, distinct SPT phases can be systematically classified by group cohomology of $G$, while for free fermions, the SPT phases can be systematically described by K-theory or homotopy group theory, which leads to the well known periodic table for topological insulators and superconductors.

Most 3D topological insulators have to be protected by some other symmetries, such as time reversal, particle hole or chiral symmetry, and the $U(1)$ charge conservation symmetry. A peculiar exception occurs when the Hamiltonian has just two effective bands. In this case, interesting topological phases, the so-called Hopf insulators, may exist. These Hopf insulator phases have no symmetry other than the prerequisite $U(1)$ charge conservation. To elucidate why this happens, let us consider a generic band Hamiltonian in 3D with $m$ filled bands and $n$ empty bands. Without symmetry constraint, the space of such Hamiltonians is topologically equivalent to the Grassmannian manifold $G_{m,m+n}$ and can be classified by the homotopy group of this Grassmannian. Since the homotopy group $\pi_3(G_{m,m+n}) = \{0\}$ for all $(m,n) \neq (1,1)$, there exists no nontrivial topological phase in general. However, when $m = n = 1$, $G_{1,2}$ is topologically equivalent to $S^2$ and the well-known Hopf map in mathematics shows that $\pi_3(G_{1,2}) = \pi_3(S^2) = \mathbb{Z}^2$. This explains why the Hopf insulators may exist only for Hamiltonians with two effective bands. The classification theory shows that the peculiar Hopf insulators may exist in 3D, but it does not tell us which Hamiltonian can realize such phases. It is even a valid question whether these phases can appear at all in physically relevant Hamiltonians. Moore, Ran, and Wen made a significant advance in this direction by constructing a Hamiltonian that realizes a special Hopf insulator with the Hopf index $\chi = 1$.

In this Rapid Communication, we construct a class of tight-binding Hamiltonians that realize arbitrary Hopf insulator phases with any integer Hopf index $\chi$. The Hamiltonians depend on two parameters and contain spin-dependent and spin-flip hopping terms. We map out the complete phase diagram and show that all the Hopf insulators can be realized with this type of Hamiltonian. We numerically calculate the surface states for these Hamiltonians and show that they have zero energy modes that are topologically protected and robust to arbitrary random perturbations with no other than the $U(1)$ symmetry constraint.

To begin with, let us notice that any two-band Hamiltonian in 3D with one filled band can be expanded in the momentum space with three Pauli matrices $\sigma = (\sigma^x, \sigma^y, \sigma^z)$ as

$$\mathcal{H}(k) = \mathbf{u}(k) \cdot \sigma,$$

where we have ignored the trivial energy-shifting term $u_0(k)I_2$ with $I_2$ being the $2 \times 2$ identity matrix. By diagonalizing $\mathcal{H}(k)$, we have the energy dispersion $E(k) = \pm |\mathbf{u}(k)|$, where $|\mathbf{u}(k)| = \sqrt{u_2^2(k) + u_3^2(k) + u_1^2(k)}$. The Hamiltonian is gapped if $|\mathbf{u}(k)| > 0$ for all $k$. For the convenience of discussion of topological properties, we denote $\mathbf{u}(k) = |\mathbf{u}(k)| (x(k), y(k), z(k))$ with $x^2(k) + y^2(k) + z^2(k) = 1$. Topologically, the Hamiltonian (1)
can be considered as a map from the momentum space \( k = (k_x, k_y, k_z) \) characterized by the Brillouin zone \( T^3 \) (\( T \) denotes a circle and \( T^3 \) is the 3D torus) to the parameter space \( u(k) \propto (x(k), y(k), z(k)) \) characterized by the Grassmannian \( G_{2,1,2} = S^2 \). Topologically distinct band insulators correspond to different classes of maps from \( T^3 \to S^2 \).

The classification of all the maps from \( T^3 \to S^2 \) is related to the torus homotopy group \( \tau_3(S^2) \). To construct non-trivial maps from \( T^3 \to S^2 \), we take two steps, first from \( S^1 \to S^2 \) and then from \( T^3 \to S^1 \). We make use of the following generalized Hopf map \( f : S^3 \to S^2 \) known in the mathematical literature:

\[
x + iy = 2\lambda \eta_{x}^{2} \eta_{y}^{2}, \quad z = \lambda(|\eta_{x}|^{2}p - |\eta_{y}|^{2}q),
\]

where \( p, q \) are integers prime to each other and \( \eta_{x}, \eta_{y} \) are complex coordinates for \( \mathbb{R}^4 \) satisfying \( |\eta_{x}|^2 + |\eta_{y}|^2 = 1 \) with the normalization \( \lambda = 1/(|\eta_{x}|^{2p} + |\eta_{y}|^{2q}) \). Equation (2) maps the coordinates \((\text{Re}[\eta_{x}], \text{Im}[\eta_{x}], \text{Re}[\eta_{y}], \text{Im}[\eta_{y}])\) of \( S^3 \) to the coordinates \((x, y, z)\) of \( S^2 \) with \( x^2 + y^2 + z^2 = 1 \). The Hopf index for the map \( f \) is known to be \( \pm pq \) with the sign determined by the orientation of \( S^3 \). We then construct another map \( g : T^3 \to S^3 \) (up to a normalization), defined by the equation

\[
\eta_{x}(k) = \sin k_{x} + it \sin k_{y}, \quad \eta_{y}(k) = \sin k_{z} + i(\cos k_{x} + \cos k_{y} + \cos k_{z} + h),
\]

where \( t \) and \( h \) are constant parameters. The composite map \( f \circ g \) from \( T^3 \to S^2 \) then defines the parameters \( u(k) \propto (x(k), y(k), z(k)) \) in the Hamiltonian as a function of the momentum \( k \). From Eqs. (2) and (3), we have \( u(k) = |u(k)|(x(k), y(k), z(k)) = (\text{Re}[2\eta_{x}^{2} \eta_{y}^{2}], \text{Im}[2\eta_{x}^{2} \eta_{y}^{2}], |\eta_{x}|^{2p} - |\eta_{y}|^{2q}) \), with \( |u(k)| = 1/\lambda(k) \). The Hamiltonian \( \hat{H}(k) = u(k) \cdot \sigma \) is \((p+q)\)th order polynomials of \( \sin(k) \) and \( \cos(k) \), which corresponds to a tight-binding model when expressed in the real space. The Hamiltonian contains spin-orbital coupling with spin-dependent hopping terms. When we choose \( p = q = 1 \) and \((t, h) = (1, -3/2)\), the Hamiltonian (1) reduces to the special case studied in Ref.\(^\text{13}\).

When the Hamiltonian is gapped with \( |u(k)| > 0 \), one can define a direction on the unit sphere \( u(k) = (u_{x}(k), u_{y}(k), u_{z}(k))/|u(k)| = (x(k), y(k), z(k)) \). From \( u(k) \), we define the Berry curvature \( F_{\mu} = \frac{i}{\hbar} \epsilon_{\mu\nu\rho} \hat{u} \cdot (\partial_{\nu} \hat{u} \times \partial_{\rho} \hat{u}) \), where \( \epsilon_{\mu\nu\rho} \) is the Levi-Civita symbol and a summation over the same indices is implied. A 3D torus \( T^3 \) has three orthogonal cross sections perpendicular to the axis \( x, y, z \), respectively. For each cross section of space \( T^2 \), one can introduce a Chern number \( C_{\mu} = \int_{T^{2}} \frac{1}{2\pi} F_{\mu} \cdot dk_{\nu}dk_{\lambda}F_{\nu\lambda} \), where \( \mu = x, y, z \) and \( \rho, \lambda \) denote directions orthogonal to \( \mu \). To classify the maps from \( T^3 \to S^2 \) represented by \( u(k) \), a topological index, the so-called Hopf index, was introduced by Pontryagin\(^\text{16}\), who showed that the Hopf index takes values in the finite group \( \mathbb{Z}_{\mathbb{Z}_{2}}: \text{GCD}(C_{x}, C_{y}, C_{z}) \) when the Chern numbers \( C_{\mu} \) are nonzero\(^\text{16}\), where \( \text{GCD} \) denotes the greatest common divisor. If the Chern numbers \( C_{\mu} = 0 \) in all three directions, the Hopf index takes all integer values \( \mathbb{Z} \) and has a simple integral expression\(^\text{15,17}\):

\[
\chi(\hat{u}) = -\int_{BZ} F \cdot dk,
\]

where \( A \) is the Berry connection (or called the gauge field) which satisfies \( \nabla \times A = F \). The Hopf index \( \chi(\hat{u}) \) is gauge invariant although its expression depends on \( A \). As we will analytically prove in the Appendix, the Chern numbers \( C_{\mu} = 0 \) for the map \( u(k) \) defined above in this paper in the gapped phase, so we can use the integral expression of Eq. (4) to calculate the Hopf index \( \chi(\hat{u}) \). The index \( \chi(\hat{u}) \) can be calculated numerically through discretization of the torus \( T^3 \). Using this method, we have numerically computed the Hopf index \( \chi(\hat{u}) \) for the Hamiltonian \( \hat{H}(k) \) with various \( p \) and \( q \), and the results are shown in Fig. 1. As the grid number increases in discretization, we see that the Chern numbers quickly drop to zero and the Hopf index approaches the integer values \( \pm pq \) or \( \pm 2pq \) depending on the parameters \( t, h \).

Based on the numerical results of \( \chi(\hat{u}) \), we construct the phase diagrams of the Hamiltonian (1) for various \( p, q \) in Fig. 2. The phase boundaries are determined from the gapless condition. The phase diagrams exhibit regular patterns: they are mirror symmetric with respect to the axis \( h = 0 \) and anti-symmetric with respect to the axis \( t = 0 \). When \( |h| > 3 \), we only have a topologically trivial phase with \( \chi(\hat{u}) = 0 \). From the result, we see that \( \chi(\hat{u}) \) has an analytic expression with \( \chi(\hat{u}) = \pm pq \) when \( 1 < |h| < 3 \) and \( \chi(\hat{u}) = \pm 2pq \) when \( |h| < 1 \).

To understand this result, we note that \( u(k) \) is a composition of two maps \( f \circ g(k) \). The generalized Hopf maps \( f \) from \( S^3 \to S^2 \) has a known Hopf index \( \pm pq \). The maps \( g \) from \( T^3 \to S^3 \) can be classified by the torus homotopy group \( \tau_3(S^3) \) and a topological invariant has been introduced to describe this classification\(^\text{18}\).
which has an integral expression
\[ \Gamma(g) = \frac{1}{12\pi^2} \int_{BZ} d\mathbf{k} \epsilon_{\alpha\beta\gamma\rho} \epsilon_{\mu\nu\tau} \frac{1}{|\eta|} \eta_{\alpha} \partial_{\beta} \eta_{\gamma} \eta_{\rho} \partial_{\tau} \eta_{\nu}, \]
where \( \eta = (\text{Re}[\eta], \text{Im}[\eta], \text{Re}[\eta], \text{Im}[\eta]) \). Direct calculation of \( \Gamma(g) \) leads to the following result:
\[ \Gamma(g) = \begin{cases} 
0, & |h| > 3 \\
1, & 1 < |h| < 3 \text{ and } t > 0 \\
-2, & |h| < 1 \text{ and } t > 0.
\end{cases} \]
Consequently, we have \( \chi(\hat{u}) = \Gamma(g) \chi(f) = \pm pq \Gamma(g) \), which is exactly the result shown in the phase diagrams in Fig. 2. A geometric interpretation is that \( \Gamma(g) \) counts how many times \( T^3 \) wraps around \( S^1 \) under the map \( g \), and \( \chi(f) \) describes how many times \( S^3 \) wraps around \( S^2 \) under the generalized Hopf map \( f \). Their composition gives the Hopf index \( \chi(\hat{u}) \). A sign flip of \( t \) changes the orientation of the sphere \( S^3 \), which induces a sign flip in \( \chi(\hat{u}) \) and produces the anti-symmetric phase diagram with respect to the axis \( t = 0 \). As \( (p, q) \) are arbitrary coprime integers, \( \chi(\hat{u}) \) apparently can take any integer value depending on the values of \( p, q \) and \( t, h \). As a consequence, the Hamiltonian \( H_k \) constructed in this communication can realize arbitrary Hopf insulator phases.

The nontrivial topological invariant guarantees existence of gapless surface states at a smooth (i.e., adiabatic) boundary between a Hopf insulator and a trivial insulator (or vacuum). Numerically, we find that gapless surface states are still present even for sharp boundaries, although we do not have an intuitive explanation why this is necessarily so as the number of bands is not well-defined at a sharp boundary and the two-band condition required for existence of the Hopf insulator could be violated at the surface. Our results are summarized in Fig. 3. From the figure, surface states and localized zero-energy modes are prominent. These surface states are topologically protected and robust under arbitrary random perturbations that only respect the prerequisite \( U(1) \) symmetry. This can be clearly seen from

FIG. 2. (Color online) Phase diagrams of the Hamiltonian for different \((p, q)\). The values of \((p, q)\) in (a), (b), (c), and (d) are chosen to be \((1, 1), (1, 2), (3, 1), \) and \((2, 3)\), respectively.

FIG. 3. (Color online) Surface states and zero-energy modes in the (001) direction for a 200-site-thick slab. The parameters \( t \) and \( h \) are chosen as \((t, h) = (1, 1.5)\) for all the figures. We have \((p, q) = (1, 2)\) for (a,b) and \((p, q) = (1, 3)\) for (c,d). In Fig. (b,d), we add random perturbations to the Hamiltonian, but otherwise keep the same parameters as (a,c). The left diagrams in (a,b,c,d) plot the energy spectrum of all 400 states at a fixed \((k_x, k_y) = (0.72, 0.72)\) for easy visualization. The points inside the gap represent the energies of the surface states. There are four (six) surface states in (a,b) ((c,d)), respectively. The right diagrams in in (a,b,c,d) show the wave functions of a surface state (upper one) and a bulk state (lower one).
Fig. 3: while the wave functions of the bulk states change dramatically under random perturbations, the wave functions of the surface states remain stable and are always sharply peaked at the boundary. This verifies that the Hopf insulators are indeed 3D topological phases. Besides the results shown in Fig. 3, we have calculated the surface states for a number of different choices of parameters \((p, q)\) and \((t, h)\), and the results consistently demonstrate that the surface states and zero energy modes are always present and robust even to substantial perturbations unless the bulk gap closes. Moreover, we roughly have more surfaces states when the absolute value of the walking distance and direction in tum walks becomes larger. However, this is not always true. A direct correspondence between the Hopf index and the total winding number of surface states may exist and deserves to be further investigated. It is also worthwhile to mention that these surface states are extended/metallic in a clean crystal, as discussed in Ref. 13, but how disorder will affect these states is an important topic that deserves further studies. The surface states might not be metallic with disorder since there is no obvious way to protect these surface state from localization without adding symmetries such as time-reversal.

An important and intriguing question is how to realize these Hopf insulators in experiments. Laser assisted hopping of ultracold atoms in an optical lattice offers a powerful tool to engineer various kinds of spin-dependent tunneling terms, and thus provides a good candidate for their realizations although the details still need to be worked out. Dipole interaction between polar molecules in optical lattices also offers possibilities to realize effective spin-dependent hopping. As argued in Ref. 13, frustrated magnetic compounds such as \(X_2Mo_2O_7\) with \(X\) being a rare earth ion are other potential candidates. In addition, Hopf insulators may be realized in 3D quantum walks, where various hopping terms are implemented by varying the walking distance and direction in each spin-dependent translation and the robust surface states can be observed with split-step schemes.

In conclusion, we have introduced a class of tight-binding Hamiltonians that realize arbitrary Hopf insulators. The topologically protected surface states and zero-energy modes in these exotic phases are robust to random perturbations that only respect the \(U(1)\) charge conservation symmetry. They are 3D topological phases and sit outside of the periodic table for topological insulators and superconductors.

**Appendix.** Here, we prove that the Chern numbers \(C_n = 0\) in all three directions for our Hamiltonian. Let us first consider \(C_x\). To prove \(C_x = \int_{-\pi}^\pi \int_{-\pi}^\pi dk_y dk_z F_x(k_y, k_z) = 0\), it is sufficient to show \(F_x(k_y, k_z) = -F_x(-k_y, -k_z)\), i.e., the function \(F_x\) has an odd parity under the exchange \((k_y, k_z) \rightarrow (-k_y, -k_z)\). We denote the parity of a given function \(F(k_y, k_z)\) as \(P[F] = \{1, -1\}\) corresponding to \{even, odd\} parity. Our aim is to prove \(P[F_x] = -1\). We let \(g_1 = Re(\eta_1(k)) = sin k_z\), \(g_2 = Im(\eta_1(k)) = t sin k_y\), \(g_3 = Re(\eta_1(k)) = sin k_z\), and \(g_4 = Im(\eta_1(k)) = cos k_x + cos k_y + cos k_z + h\). Apparently, \(P[g_1] = P[g_4] = 1\) and \(P[g_2] = P[g_3] = -1\).

We can normalize the \(g\)-vector as \(\hat{g} = g/|g| = (g_1, g_2, g_3, g_4)/\sqrt{g_1^2 + g_2^2 + g_3^2 + g_4^2}\). The components of \(\hat{g}\) have the same parity as the unnormalized ones. From the definition, we have \(\hat{u}_x = \text{Re}\left[2\hat{\lambda}(g_1 + ig_2)^2(g_3 - ig_4)^q\right] = 2\lambda \text{Re}\left[\sum_{\alpha=0}^p \sum_{\beta=0}^q C_{\alpha\beta}^p q^{\beta}(1)^{p+q-\alpha-\beta} g_1^\alpha g_2^{p-q-\alpha-\beta} g_3^\alpha g_4^{p-q-\alpha-\beta}\right]\), where \(C_{\alpha\beta}^p\) (\(C_{\beta\alpha}^p\)) denote the binomial coefficients and \(\lambda = 1/((\hat{g}_1 + i\hat{g}_2)^2 + (\hat{g}_3 + i\hat{g}_4)^2)\). The exponent \(p+q-\alpha-\beta\) of \(i\) in \(\hat{u}_x\) has to be even to have a nonzero real part, so \(P[\hat{u}_x] = P[g_2^p g_3^q g_4^{p-q}] = (-1)^q\). Similarly, by using \(\hat{u}_y = \text{Im}[2\hat{\lambda}(g_1 + ig_2)^2(g_3 - ig_4)^q]\), we find \(P[\hat{u}_y] = -P[\hat{u}_x]\). Finally, from \(\hat{u}_z = \hat{\lambda}((\hat{g}_1 + i\hat{g}_2)^2 - (\hat{g}_3 + i\hat{g}_4)^2)\) we obtain \(P[\hat{u}_z] = 1\). As a consequence, \(P[\hat{u} \cdot (\partial_x \hat{u} \times \partial_y \hat{u})] = -1\). Therefore, \(P[F_x] = P[\hat{u} \cdot (\partial_x \hat{u} \times \partial_y \hat{u})] = -1\). This proves that \(C_x = 0\). By the same parity arguments, we can show \(C_y = C_z = 0\).

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SUPPLEMENTAL MATERIAL: HOPF INSULATORS AND THEIR TOPOLOGICALLY PROTECTED SURFACE STATES

In this supplemental material, we explain the details on how to obtain the surface states and the zero energy modes.

We give more details on how to numerically calculate the surface states and zero energy modes. We take a slab in the (001) direction and maintain the periodic boundary condition in the \((x, y)\)-directions. Along the \(z\)-direction, we work in the real space by an inverse Fourier transform of the momentum \(k_z\). Suppose we consider a \(N_z\)-site-thick slab, for any fixed \((k_x, k_y)\), we arrange the \(2N_z\) basis-vectors of the Hilbert space by \((|↓⟩_1, |↑⟩_1, \cdots, |↓⟩_{N_z}, |↑⟩_{N_z})\), where the subscript denotes the site number. After the inverse Fourier transform, the Hamiltonian can be written in general as 

\[
\mathcal{H} = \sum_{k_z, k_y} \mathcal{H}_{k_z, k_y},
\]

where

\[
\mathcal{H}_{k_z, k_y} = \sum_{i=1}^{2N_z} \sum_{j=1}^{2N_z} t_{ij}^{k_z, k_y} c_{k_z, k_y, i}^\dagger c_{k_z, k_y, j}.
\]

We aim to find an analytical expression for \(t_{ij}^{k_z, k_y}\). From the text, the Hamiltonian in the momentum space reads

\[
\mathcal{H} = \sum_k \Psi^\dagger(k) \mathcal{H}(k) \Psi(k) = \sum_k \{u_x c_{k, \uparrow}^\dagger c_{k, \uparrow} - u_x c_{k, \downarrow}^\dagger c_{k, \downarrow} + [(u_x + iu_y) c_{k, \downarrow}^\dagger c_{k, \uparrow} + h.c.]\},
\]

where

\[
\begin{align*}
u_x + iu_y &= 2(\sin k_x + it \sin k_y)^p[\sin k_z - i(\cos k_x + \cos k_y + \cos k_z + h)]^q \\
u_z &= (\sin^2 k_x + t^2 \sin^2 k_y)^p - \langle \sin^2 k_z + \cos k_x + \cos k_y + \cos k_z + h \rangle^2.
\end{align*}
\]

Since we only perform inverse Fourier transform in the \(z\) direction and keep \((k_x, k_y)\) in the momentum space, we can take \(k_x\) and \(k_y\) as constants. Let \(A = 2(\sin k_x + it \sin k_y)^p\) and \(B = -i(\cos k_x + \cos k_y + h)\), then Eq. (7) reduces to

\[
u_x + iu_y = A(-ie^{ik_z} + B)^q = A \sum_{\kappa=0}^q \left(\frac{q}{\kappa}\right) (-i)^\kappa e^{ik_z} B^{q-\kappa}
\]

\[
= \sum_{\kappa=0}^q D_\kappa e^{ik_z},
\]

where \(D_\kappa = A \left(\frac{q}{\kappa}\right) (-i)^\kappa B^{q-\kappa}\). Similarly, for the \(u_z\) term, we define \(R = (\sin^2 k_x + t^2 \sin^2 k_y)^p\), \(S = 1 + (\cos k_x + \cos k_y + h)^2\), \(T = (\cos k_x + \cos k_y + h)\) and \(Q = S/T\). Eq. (8) then reduces to

\[
u_z = R - T^q(Q + e^{ik_z} + e^{-ik_z})^q
\]

\[
= R - \sum_{\alpha + \beta + \kappa = q} \left(\frac{q}{\alpha, \beta, \kappa}\right) T^q Q^\kappa e^{i(\alpha-\beta)k_z}
\]

\[
= R - \sum_{\alpha + \beta + \kappa = q} J_{\alpha\beta\kappa} e^{i(\alpha-\beta)k_z},
\]

where \(\left(\frac{q}{\alpha, \beta, \kappa}\right) = \frac{q!}{\alpha! \beta! \kappa!}\) is the trinomial coefficient and \(J_{\alpha\beta\kappa} = \left(\frac{q}{\alpha, \beta, \kappa}\right) T^q Q^\kappa\). Now we are ready to perform the inverse Fourier transform in the \(z\) direction:

\[
c_{k_z, k_y, k_z, \sigma} = \frac{1}{\sqrt{N_z}} \sum_{z} e^{izk_z} c_{k_z, k_y, z, \sigma},
\]

\[
c_{k_z, k_y, k_z, \sigma}^\dagger = \frac{1}{\sqrt{N_z}} \sum_{z} e^{-izk_z} c_{k_z, k_y, z, \sigma}^\dagger.
\]
After the transformation, we obtain

\[
\mathcal{H}^{k_x, k_y} = [(Rc_{k_x, k_y, z, \uparrow}^\dagger C_{k_x, k_y, z, \uparrow} - \sum_{\alpha+\beta+\kappa=q} J_{\alpha\beta\kappa \eta} c_{k_x, k_y, z, \uparrow}^\dagger C_{k_x, k_y, z-\alpha-\beta, \uparrow}) - (\uparrow\rightarrow\downarrow)]
\]

\[
+ \left[ \sum_{\kappa=0}^q D_{\kappa} c_{k_x, k_y, z, \downarrow}^\dagger C_{k_x, k_y, z-\kappa, \uparrow} + \text{h.c.} \right].
\]

Comparing Eq. (10) with Eq. (5), we find the expressions

\[
l^{k_x, k_y}_{2k, 2l} = R\delta_{k, l} - \sum_{\alpha+\beta+\kappa=q} J_{\alpha\beta\kappa \eta} \delta_{2k, 2l+\alpha-\beta},
\]

\[
l^{k_x, k_y}_{2k-1, 2l-1} = -R\delta_{k, l} + \sum_{\alpha+\beta+\kappa=q} J_{\alpha\beta\kappa \eta} \delta_{2k, 2l+\alpha-\beta},
\]

\[
l^{k_x, k_y}_{2k-1, 2l} = \sum_{\kappa=0}^q D_{\kappa} \delta_{2k-1, 2l+\kappa},
\]

\[
l^{k_x, k_y}_{2k, 2l-1} = \sum_{\kappa=0}^q D_{\kappa}^* \delta_{2k+1, 2l-1},
\]

where \(0 \leq k, l \leq N_z\). Hence, for each \(k_x\) and \(k_y\), we have a \(2N_z \times 2N_z\) matrix \(t^{k_x, k_y}\) with its \((i, j)\)-th entry \(t^{k_x, k_y}_{ij}\).

Numerically diagonalizing this matrix for fixed \(k_x\) and \(k_y\), we obtain the energy spectrum of \(2N_z\) states. For each \(k_x\) and \(k_y\), we count the number of surface states by noticing that surface state energies have huge gaps from the bulk state energies. Fig. 4 shows the number of edge states for all \(k_x\) and \(k_y\) values by imposing a minimum relative separation from the bulk. The number of surface states is counted as the union of all edge states for all \((k_x, k_y)\). The right diagram shows the case where small random perturbations are included. We see that the surface states are robust to random perturbations without any symmetry constraint. In the text, for easy visualization, we plotted the energy spectrum of the Hamiltonian in Fig. 3 at fixed \((k_x, k_y) \approx (0.72, 0.72)\). Each in-gap point corresponds to a surface state.