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Phys. Rev. B **96**, 041103 — Published 5 July 2017

DOI: [10.1103/PhysRevB.96.041103](https://doi.org/10.1103/PhysRevB.96.041103)

Nodal-link semimetals

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In topological semimetals, the valance band and conduction band meet at zero-dimensional nodal points or one-dimensional nodal rings, which are protected by band topology and symmetries. In this Rapid Communication, we introduce “nodal-link semimetals”, which host linked nodal rings in the Brillouin zone. We put forward a general recipe based on the Hopf map for constructing models of nodal-link semimetal. The consequences of nodal ring linking in the Landau levels and Floquet properties are investigated.

Topological phases of matter have been among the most active research subjects in condensed matter physics. They can be broadly classified as two major classes. The first class of phases, including topological insulators and superconductors[1–6], and other symmetry protected topological phases[7], have gapped bulks with nontrivial topological structures characterized by topological invariants[8–15], dictating the existence of robust gapless modes on the boundary.

More recently, the second major class of topological materials, known as topological semimetals, have attracted widespread attentions. In the noninteracting limit, they are characterized by topologically robust \mathbf{k} -space band-touching manifolds, which can be zero-dimensional (0D) nodal points or one-dimensional (1D) nodal rings (or nodal lines). The bulk Dirac[16–25] and Weyl points[26–40] are responsible for novel phenomena related to chiral anomaly[41–51]. Moreover, bulk Weyl points entail surface Fermi arcs, while nodal rings[52–72] imply flat surface band (drumhead states) that may trigger interesting correlation effects[73]. Nodal rings have been predicted (e.g., Cu₃PdN[60, 61], Hg₃As₂[74], Ca₃P₂[63, 75], 3D carbon networks[58], CaP₃[76], Alkali Earth Metals[77, 78]) and experimentally studied in quite a few materials (e.g., PbTaSe₂[62], ZrSiTe[79], ZrSiS[80–84]). Notably, nodal rings can be driven to Floquet Weyl points by circularly polarized light[85–89], accordingly, the drumhead surface states become Fermi arcs.

Unlike nodal points, nodal rings allow richer topological structures. They can touch at special points[60, 61, 90, 91], enabling formations of nodal chains[92, 93]. In this Letter, we introduce new types of topological semimetals, dubbed “nodal-link semimetals”, which host nontrivially linked nodal rings[e.g., Fig.1(e)]. Furthermore, a method is introduced for constructing two-band models of nodal-link semimetals. We investigate generic physical consequences of nontrivial linking; in particular, a global toroidal π Berry phase generates a half-integer shift of Landau level index when the magnetic field is perpendicular to the ring plane. In addition, a suitable periodic external field can drive nodal-link semimetal to a Floquet Hopf insulator.

Models.—Nodal rings come from the crossing of two adjacent bands, thus we focus on two-band Bloch Hamiltonians,

which can generally be written as

$$H(\mathbf{k}) = a_0(\mathbf{k})\mathbf{1} + a_1(\mathbf{k})\tau_x + a_2(\mathbf{k})\tau_y + a_3(\mathbf{k})\tau_z, \quad (1)$$

where $\mathbf{k} = (k_x, k_y, k_z)$, τ_i 's are Pauli matrices, and $a_0(\mathbf{k}) = 0$ will be adopted for simplicity (nonzero a_0 can be trivially included, if needed). Nodal rings are protected by crystal symmetries. For concreteness, we take the PT symmetry[66, 94] that ensures the reality of $H(\mathbf{k})$, i.e., $a_2(\mathbf{k}) = 0$. Now the spectra are $E_{\pm}(\mathbf{k}) = \pm \sqrt{a_1^2 + a_3^2}$, and the nodal rings are given by solving $a_1(\mathbf{k}) = a_3(\mathbf{k}) = 0$. A purpose of this Letter is to construct models with mutually linked nodal rings.

Instead of taking trial-and-error approaches, we put forward a general method based on Hopf maps[95, 96]. They play special roles in quantum spin systems[95, 97], topological Hopf insulators[98–103], liquid-crystal solitons[104], quench dynamics of Chern insulators[105], and minimal models for topologically trivial superconductor-based Majorana zero modes[106]. Mathematically, a Hopf map is a nontrivial mapping from a 3-sphere S^3 to a 2-sphere S^2 , which possesses a nonzero Hopf invariant[95, 96, 98]. Moreover, it has the geometrical property that the preimage circles of any two points on S^2 are linked. Mappings from a 3D torus T^3 to S^2 inherit the nontrivial topology of Hopf maps through $T^3 \rightarrow S^3 \rightarrow S^2$, where $T^3 \rightarrow S^3$ is a map with unit winding number, and $S^3 \rightarrow S^2$ is a Hopf map.

Given any vector function $\mathbf{d}(\mathbf{k}) = (d_x, d_y, d_z)$ on the Brillouin zone T^3 , one can define a mapping from T^3 to S^2 by $\mathbf{k} \rightarrow \hat{\mathbf{d}}(\mathbf{k})$, where $\hat{\mathbf{d}} \equiv \mathbf{d}/|\mathbf{d}|$. To define the Hopf invariant, it is convenient to express the vector \mathbf{d} in terms of a spinor, namely, $d_i(\mathbf{k}) = z^\dagger \tau_i z$, with $z(\mathbf{k}) = (z_1, z_2)^T$. Let us write $z_1 = N_1 + iN_2, z_2 = N_3 + iN_4$, then the Hopf invariant simplifies to[103]

$$n_h = \frac{1}{2\pi^2} \int d^3k \epsilon^{abcd} \hat{N}_a \partial_{k_x} \hat{N}_b \partial_{k_y} \hat{N}_c \partial_{k_z} \hat{N}_d, \quad (2)$$

where \hat{N}_a is the a -th component of the vector $\mathbf{N} = (N_1, N_2, N_3, N_4)$ normalized to unit length.

Let us take a point on S^2 , say $\hat{\mathbf{n}}_1 = (0, 1, 0)$. Under the mapping $\mathbf{k} \rightarrow \hat{\mathbf{d}}(\mathbf{k})$, all the preimages of $\hat{\mathbf{n}}_1$ have to satisfy

$$d_x(\mathbf{k}) = d_z(\mathbf{k}) = 0, \quad (3)$$

however, the converse is not true, namely, a solution of Eq.(3) is not necessarily a preimage of $\hat{\mathbf{n}}_1$. In fact, the preimages

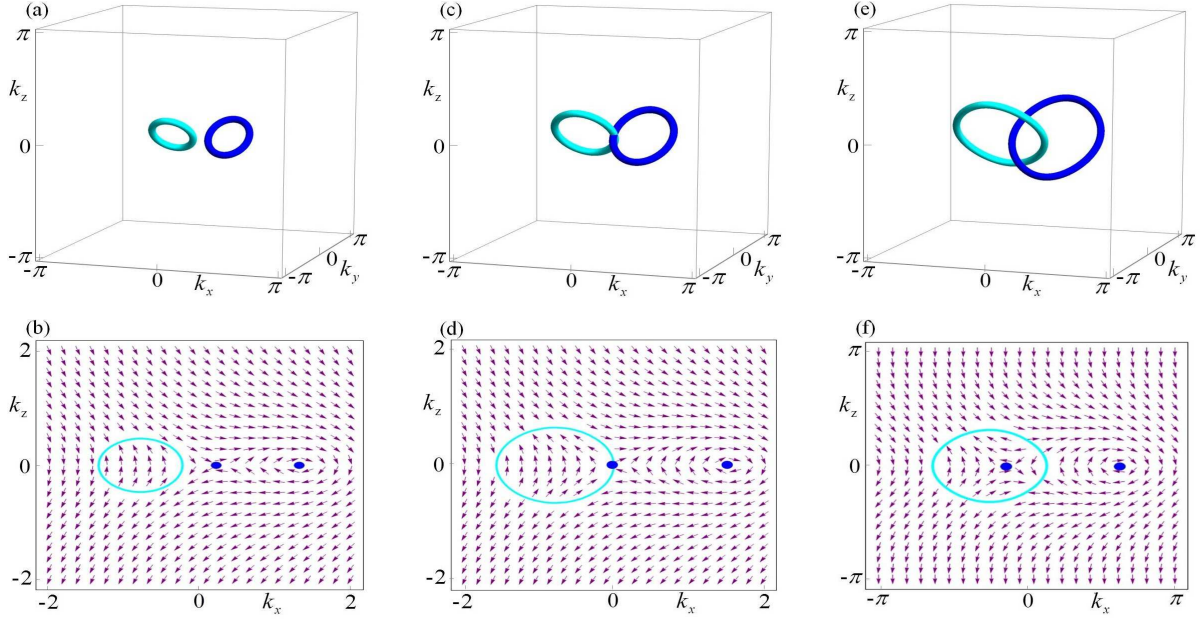


FIG. 1. Nodal rings of Eq.(6), and the pseudospin textures in the $k_y = k_z$ plane, for (a,b) $m_0 = 3.2$, (c,d) $m_0 = 3.0$, and (e,f) $m_0 = 2.5$. The light blue ring locates in the $k_y = k_z$ plane, while the dark blue ring is perpendicular to it.

of $\hat{\mathbf{n}}_2 = (0, -1, 0)$ also satisfy Eq.(3). In a single equation, Eq.(3) gives the preimages of both $\hat{\mathbf{n}}_1$ and $\hat{\mathbf{n}}_2$. This is among key observations in our construction. When n_h is nonzero, the preimages circles of any two points (say $\hat{\mathbf{n}}_1$ and $\hat{\mathbf{n}}_2$) are linked. Therefore, we can obtain linked nodal rings by taking

$$a_1(\mathbf{k}) = d_x(\mathbf{k}), \quad a_3(\mathbf{k}) = d_z(\mathbf{k}). \quad (4)$$

Recall that the a_2 term is absent due to crystal symmetries, as discussed above. The same method also works if we start from a different $\mathbf{n}_{1,2}$; for instance, taking $\mathbf{n}_{1,2} = (0, 0, \pm 1)$ leads to $a_1 = d_x, a_3 = d_y$, which yields a nodal link as well.

This is a quite general approach to construct nodal-link semimetals. As an example, we take[98–103]

$$\begin{aligned} N_1 &= \sin k_x, \quad N_2 = \sin k_y, \quad N_3 = \sin k_z, \\ N_4 &= \cos k_x + \cos k_y + \cos k_z - m_0, \end{aligned} \quad (5)$$

which has $n_h = -1$ for $1 < m_0 < 3$ and $n_h = 0$ for $m_0 > 3$. The explicit forms of d_i 's read

$$\begin{aligned} d_x &= 2 \sin k_x \sin k_z + 2 \sin k_y \left(\sum_{i=x,y,z} \cos k_i - m_0 \right), \\ d_y &= -2 \sin k_y \sin k_z + 2 \sin k_x \left(\sum_{i=x,y,z} \cos k_i - m_0 \right), \\ d_z &= \sin^2 k_x + \sin^2 k_y - \sin^2 k_z - \left(\sum_{i=x,y,z} \cos k_i - m_0 \right)^2. \end{aligned}$$

Following Eq.(4), a lattice model of nodal-link semimetal is

$$\begin{aligned} H(\mathbf{k}) &= [2 \sin k_x \sin k_z + 2 \sin k_y \left(\sum_{i=x,y,z} \cos k_i - m_0 \right)] \tau_x \\ &+ [\sin^2 k_x + \sin^2 k_y - \sin^2 k_z - \left(\sum_{i=x,y,z} \cos k_i - m_0 \right)^2] \tau_z. \end{aligned} \quad (6)$$

The nodal rings consist of points where both coefficients of τ_x and τ_z vanish. We find that one of the rings is $k_y = k_z, \sin k_x = m_0 - \sum_i \cos k_i$, and the other is $k_y = -k_z, \sin k_x = \sum_i \cos k_i - m_0$, shown in light and dark blue, respectively, in Fig.1. Fig.1(a) shows the unlinked rings for $m_0 = 3.2$, and Fig.1(e) illustrates the linked rings for $m_0 = 2.5$ (a Hopf link). The critical point $m_0 = 3.0$, when the two rings cross each other, is shown in Fig.1(c). The direction of pseudospin vector (d_x, d_z) is plotted in Fig.1(b)(d)(f), indicating that the light blue ring encloses a pseudospin vortex in the linked regime [Fig.1(f)], in contrast to the unlinked case [Fig.1(b)]. The surface states for $m_0 = 2.5$ are shown in Fig.2. The two-disk-overlapping region has zero and two flat bands in Fig.2(a) and (c), respectively, which is consistent with the winding number[107] in each region.

Near the critical point[Fig.1(c)], we can expand the Hamiltonian as:

$$\begin{aligned} H(\mathbf{k}) &= [2k_x k_z + 2k_y(m - k^2/2)] \tau_x \\ &+ [k_x^2 + k_y^2 - k_z^2 - (m - k^2/2)^2] \tau_z, \end{aligned} \quad (7)$$

where $k^2 = \sum_{i=x,y,z} k_i^2, m = 3 - m_0$. At the critical point $m = 0$, we have $H(\mathbf{k}) \approx 2k_x k_z \tau_x + (k_x^2 + k_y^2 - k_z^2) \tau_z$, thus the dispersion is quadratic in all three directions. Consequently, we find that, for $m = 0$, the density of states follows $g(E) \sim \sqrt{E}$ near zero energy, in contrast to $g(E) \sim E$ for $m \neq 0$. Just like that nontrivial topology of insulators can be undone by closing the energy gap, the nontrivial nodal-line linking can be untied through quadratic-dispersion critical points[Fig.1(c)], where $g(E) \sim \sqrt{E}$.

We remark that, although a nodal chain[92, 93] also contains crossings like Fig.1(c), its dispersion is not quadratic in all three directions at the crossing point, and the density of

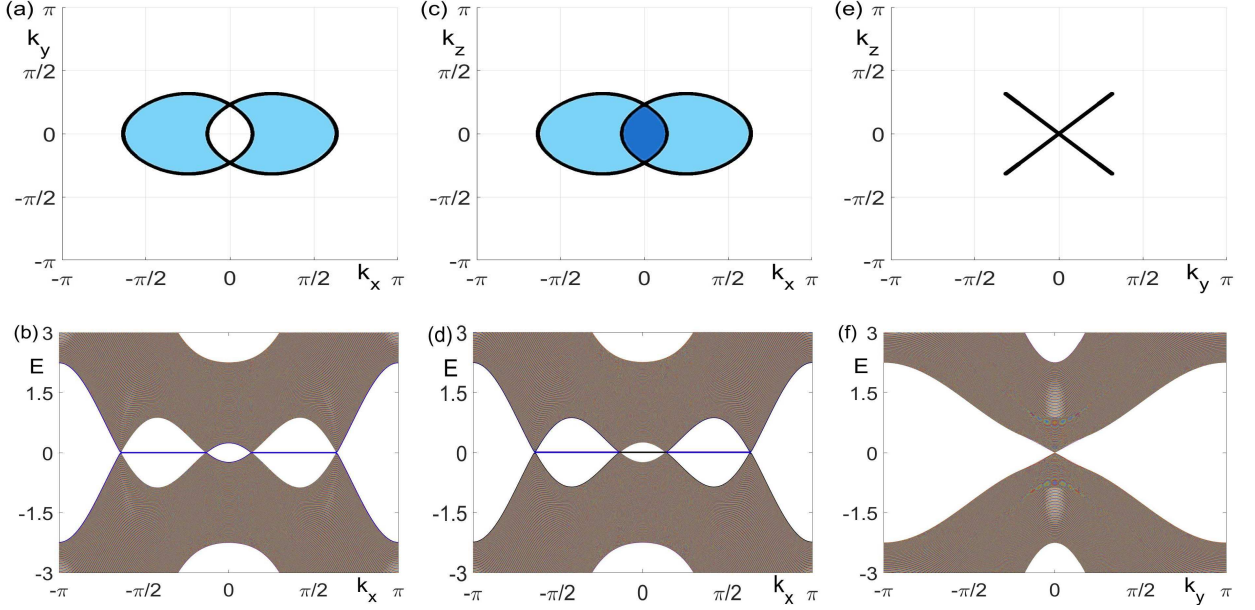


FIG. 2. (a) Surface flat bands for $z = 0$ boundary. (b) The spectra as a function of k_x (fixed $k_y = 0$) for a 800-site-thick slab perpendicular to z axis. (c) The $y = 0$ surface states. (d) Spectra of a slab perpendicular to y axis. (e) $x = 0$ surface Brillouin zone without surface states. (f) Spectra of a slab perpendicular to x axis. The black rings or lines in (a,c,e) are projections of the bulk nodal link to the surface Brillouin zone. In (c), the number of bands in the two-disk-overlapping region is twice that of the non-overlapping regions. Here, $m_0 = 2.5$.

states is linear instead of square-root.

Landau levels.—A key difference between a usual unlinked ring and a linked one is a global Berry phase along the ring. Let us draw a thin torus enclosing a nodal ring, then the Berry phase along the poloidal direction is always π ; in contrast, the Berry phase along the toroidal direction can be 0 or π (mod 2π), corresponding to the unlinked and linked ring, respectively [The 0 or π toroidal Berry phase of the light-blue ring can be read from the spin texture in Fig.1(b) and (f), respectively].

This π toroidal Berry phase can qualitatively affect the Landau levels. It is challenging to find analytic expressions of Landau levels for Eq.(6), nevertheless, a different model allows a simple solution: we take $a_1 = d_x(\mathbf{k})$, $a_3 = d_y(\mathbf{k})$ in Eq.(1). This Bloch Hamiltonian harbors four straight nodal lines at $(k_x, k_y) = (0, 0)$, $(0, \pi)$, $(\pi, 0)$ and (π, π) , respectively. When $1 < m_0 < 3$, a nodal ring encircling $(k_x, k_y) = (0, 0)$ is also found as $\cos k_x + \cos k_y = m_0 - 1$ in the $k_z = 0$ plane. Instead of working on this lattice Hamiltonian, we study its continuum limit for simplicity:

$$H(\mathbf{k}) = 2[k_x k_z + (m - Ck^2)k_y]\tau_x + 2[-k_y k_z + (m - Ck^2)k_x]\tau_y, \quad (8)$$

where $k^2 = k_x^2 + k_y^2 + k_z^2$. $m = 3 - m_0$, and $C = 0.5$. We have also done a basis change $\tau_z \rightarrow \tau_y$ for later convenience. This model hosts a nodal line $k_x = k_y = 0$ and a nodal ring $k_x^2 + k_y^2 = m/C$ in the $k_z = 0$ plane, which are linked.

Now we add a magnetic field along the z direction, $\mathbf{B} = B\hat{z}$. It is standard to do the replacement $\mathbf{k} \rightarrow \mathbf{\Pi} = -i\nabla + e\mathbf{A}$ with $\mathbf{A} = (0, Bx, 0)$ [108]. It is convenient to introduce the ladder operators, $a = \frac{l_B}{\sqrt{2}}(\Pi_x - i\Pi_y)$, $a^\dagger = \frac{l_B}{\sqrt{2}}(\Pi_x + i\Pi_y)$, where $l_B =$

$1/\sqrt{eB}$ is the magnetic length. The Hamiltonian becomes

$$H = \begin{pmatrix} 0 & f(k_z) \frac{\sqrt{2}a^\dagger}{l_B} \\ \frac{\sqrt{2}a}{l_B} f^\dagger(k_z) & 0 \end{pmatrix} \quad (9)$$

where $f(k_z) = 2[k_z - i(m - Ck_z^2 - \frac{2Ca^\dagger a}{l_B^2})]$. The Landau levels are found to be

$$E_{n,\pm}(k_z) = \begin{cases} \pm \sqrt{8n[k_z^2 + (m - Ck_z^2 - n\omega_c)^2]}/l_B, & n \geq 1, \\ 0, & n = 0, \end{cases} \quad (10)$$

where $\omega_c \equiv 2C/l_B^2$. The low energy eigenvalues are around $n \sim 0$ and $n \sim m/\omega_c$, the former coming from the central nodal line at $k_x = k_y = 0$, while the latter coming from the nodal ring in the $k_z = 0$ plane. As a comparison, we also consider a model with an unlinked ring:

$$H(\mathbf{k}) = (m - Ck^2)\tau_x + k_z\tau_z. \quad (11)$$

Following the same steps, we find that the Landau levels are

$$E_{n,\pm}(k_z) = \pm \sqrt{k_z^2 + [m - Ck_z^2 - (n + \frac{1}{2})\omega_c]^2}, \quad n \geq 0. \quad (12)$$

Comparing Eq.(10) and Eq.(12), we see that the presence of a linked nodal line causes a shift of Landau level index by $1/2$, namely, $n \rightarrow n - 1/2$, which is a consequence of the π toroidal Berry phase. Such a shift can be measured by magneto-transport or magneto-optical experiments.

To highlight the effect of π Berry phase, we re-derive the Landau levels using semiclassical quantization[109–111]:

$$S(k_z) = 2\pi eB(n + 1/2 - \phi_B/2\pi), \quad (13)$$

where S is the cross-sectional area of a \mathbf{k} -space orbit, and ϕ_B is the Berry phase along the orbit. For the linked ring, we have $\phi_B = \pi$ (the toroidal Berry phase), and a semiclassical calculation (see Supplemental Material for details) yields the Landau levels in Eq.(10).

Floquet Hopf insulator from nodal link.—We will show that driving the nodal-link semimetal described by Eq.(8) creates a Floquet Hopf insulator. We consider a periodic driving generated by a circularly polarized light(CPL) propagating in z direction. The vector potential $\mathbf{A}(t) = A_0(\cos \omega t, \eta \sin \omega t, 0)$, where $\eta = 1$ and -1 stands for right-handed and left-handed CPL, respectively. Its effect is described by the minimal coupling, $H(\mathbf{k}) \rightarrow H[\mathbf{k} + e\mathbf{A}(t)]$. The full Hamiltonian is time-periodic, $H(\mathbf{k}, t + T) = H(\mathbf{k}, t)$ with $T = 2\pi/\omega$, thus, it can be expanded as $H(\mathbf{k}, t) = \sum_n \mathcal{H}_n(\mathbf{k})e^{in\omega t}$, with

$$\begin{aligned} \mathcal{H}_0(\mathbf{k}) &= 2[k_x k_z + k_y(\tilde{m}_2 - Ck^2)]\tau_x \\ &\quad + 2[-k_y k_z + k_x(\tilde{m}_2 - Ck^2)]\tau_y, \\ \mathcal{H}_{\pm 1}(\mathbf{k}) &= eA_0[k_z \mp i\eta(\tilde{m}_1 - Ck^2) - 2Ck_y(k_x \mp i\eta k_y)]\tau_x \\ &\quad + eA_0[\pm i\eta k_z + (\tilde{m}_1 - Ck^2) - 2Ck_x(k_x \mp i\eta k_y)]\tau_y, \\ \mathcal{H}_{\pm 2}(\mathbf{k}) &= Ce^2 A_0^2 [(k_y \pm i\eta k_x)\tau_x - (k_x \mp i\eta k_y)\tau_y], \end{aligned} \quad (14)$$

where $\tilde{m}_{j=1,2} = m - jCe^2 A_0^2$. In the off-resonance regime, we can use an effective time-independent Hamiltonian[112–120]:

$$\begin{aligned} H_{\text{eff}}(\mathbf{k}) &= \mathcal{H}_0 + \sum_{n \geq 1} \frac{[\mathcal{H}_n, \mathcal{H}_{-n}]}{n\omega} + O\left(\frac{1}{\omega^2}\right) \\ &= 2[k_x k_z + k_y(\tilde{m}_2 - Ck^2)]\tau_x + 2[-k_y k_z + k_x(\tilde{m}_2 - Ck^2)]\tau_y \\ &\quad + \lambda[k_z^2 + (\tilde{m}_1 - Ck^2)^2 - 2Ck_p^2(\tilde{m}_1 - Ck^2 - \frac{\gamma}{4})]\tau_z, \end{aligned} \quad (15)$$

where $\lambda = 4\eta e^2 A_0^2/\omega$, $\gamma = Ce^2 A_0^2$, and $k_p = \sqrt{k_x^2 + k_y^2}$. The energy spectrum of H_{eff} is fully gapped. For weak driving, the minima of the band are located at $k_z = 0$, $Ck_p^2 = \tilde{m}_2$, and the gap is estimated as $E_g \approx 12mC(eA_0)^4/\omega$. In the CPL approach, this gap is expected to be rather small. It can be promoted to the order of $(eA_0)^2$ by adding a small $\Delta\tau_z$ pseudo-Zeeman term (see Supplemental Material for details of calculation).

If we remove the τ_z term, Eq.(15) hosts a nodal ring and a nodal line linked together. It is readily checked that the coefficient of τ_z is positive on the line and negative on the ring (for $\eta = +1$), thus, the corresponding unit $\hat{\mathbf{d}}$ vector is $(0, 0, 1)$ and $(0, 0, -1)$, respectively. This fact suggests that Eq.(15) describes a (Floquet) Hopf insulators[98–103], or more precisely, a Floquet Hopf-Chern insulator[102], because the Chern number $C(k_z) = -2$ for arbitrary k_z , as found in our numerical calculation. In the definition of Hopf invariant[95, 98], a nonsingular global Berry potential is needed, which is impossible in the presence of nonzero Chern number, nevertheless, we can study topological surface states. For a slab perpendicular to the z direction, we can solve the differential equation $H_{\text{eff}}(k_x, k_y, -i\partial_z)\Psi_{k_x, k_y}(z) = E(k_x, k_y)\Psi_{k_x, k_y}(z)$, which gives one surface band for each surface, whose dispersion is (see Supplemental Material for calculation): $E_\alpha(\mathbf{k}) = \alpha\lambda[\tilde{m}_2/C + \gamma^2 - (1 + 3C\gamma/2)k_p^2]$, where

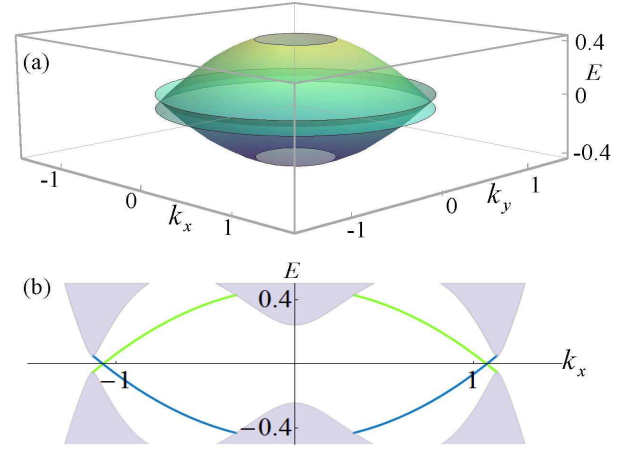


FIG. 3. Surface states of a slab perpendicular to the z axis. Parameters are $\omega = 4$, $m = 1$, $eA_0 = 0.6$, $C = 0.5$. (a) 3D view of surface bands. (b) $E(k_x)$ with $k_y = 0$ fixed. The grey regions are bulk bands.

$\alpha = + (-)$ for top (bottom) surface. Shown in Fig.3, this dispersion is characteristic of a Hopf insulator[98]. For a fixed θ (defined as $\theta = \arctan k_y/k_x$), the surface state of either top or bottom surface is chiral, which can be understood in terms of a Chern number in the (k_p, k_z) space[103].

Conclusions.—We have introduced nodal-link semimetals into the family of topological semimetals. A general method for their model construction has been put forward. These phases may be realized by tuning the hoppings in optical lattices. Finding a solid-state material will be an important progress. It will also be worthwhile to study possible novel effects of linking. Theoretically, our models lay useful ground work for topological field theories[121] in the Brillouin zone.

Note added.—After finishing this manuscript, we became aware of a related preprint by Chen et al, in which a double-helix link is constructed[122].

Acknowledgements.—We thank Chen Fang for useful comment. Z.Y., R.B., and Z.W. are supported by NSFC (No. 11674189). Z.Y. is supported in part by China Postdoctoral Science Foundation (2016M590082). L.L is supported by the Ministry of Science and Technology of China (No. 2016YFA0302400) and the National Thousand-Young-Talents Program of China. S.C.Z. is supported by NSF (No. DMR-1305677).

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