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Tsuneya Yoshida and Yasuhiro Hatsugai Phys. Rev. B **106**, 205147 — Published 28 November 2022 DOI: 10.1103/PhysRevB.106.205147

Reduction of one-dimensional non-Hermitian point-gap topology by interactions

Tsuneya Yoshida and Yasuhiro Hatsugai

Department of Physics, University of Tsukuba, Ibaraki 305-8571, Japan

(Dated: October 28, 2022)

In spite of extensive works on the non-Hermitian topology, interaction effects remain crucial questions. We hereby analyze correlated non-Hermitian systems with special emphasis on the one-dimensional point-gap topology. Specifically, our analysis elucidates that interactions result in reduction of the topological classification $\mathbb{Z} \times \mathbb{Z} \to \mathbb{Z}$ for systems of one synthetic dimension with charge U(1) symmetry and spin-parity symmetry. Furthermore, we analyze an extended Hatano-Nelson chain which exhibits striking interaction effects; interactions destroy the non-Hermitian skin effect at the non-interacting level. This fragility of the non-Hermitian skin effect against interactions is consistent with the reduction of the point-gap topology in the one spatial dimension. The above discoveries shed new light on the topology of correlated systems and open up new directions of researches on non-Hermitian topological physics.

I. INTRODUCTION

Topological insulators and superconductors have been extensively analyzed in these 15 years $^{1-9}$. In particular, considerable efforts have been devoted to understanding interaction effects on the non-trivial topology, which has revealed a variety of unique phenomena. For instance, interaction effects induce topological ordered phases^{10–16} which host anyons. In addition, it has turned out that interaction effects change Z-classification of topological superconductors at the mean-field level to \mathbb{Z}_8 classification¹⁷. Such a reduction phenomenon of possible topological phases for a given symmetry class has been theoretically reported for arbitrary spatial dimensions^{18–32}. Furthermore, a theoretical work³³ has elucidated that the reduction can also occur in synthetic dimensions which are considered to be fabricated in cold $atoms^{34-37}$. These developments reveal the ubiquity of the reduction phenomena.

Along with the above significant progress, understanding of the non-Hermitian band topology has been rapidly developed in these years^{38–43}. Remarkably, it has been elucidated that the point-gap topology induces novel phenomena which do not have Hermitian counterparts^{44–55}. A prime example is the emergence of the exceptional points^{56–60} (and their symmetry-protected variants^{61–69}) on which the point-gap topology induces band touching for both the real and the imaginary parts. Another remarkable phenomenon is a non-Hermitian skin effect which results in extreme sensitivity to the presence/absence of boundaries^{46,70–76}. So far, the non-Hermitian topological band theory has been applied to a wide range of systems from quantum^{77–88} to classical systems^{89–103}.

While most of the studies have focused on the noninteracting cases so far, interaction effects on the non-Hermitian topology attract growing interests^{104–123} due to the potential presence of novel non-Hermitian phenomena. Such interest of interaction effects on the non-Hermitian topology is further enhanced by recent development of technology in cold atoms which allows us to experimentally tune both dissipation and two-body interactions^{124,125}. Despite these efforts, current understanding of the point-gap topology in correlated systems is quite limited. In particular, the knowledge about the reduction of the point-gap topology is limited only to zero dimension¹²⁶, which poses the following significant question: fate of the higher dimensional point-gap topology under interactions.

We hereby address a primitive version of the above question. Specifically, we analyze fate of the onedimensional point-gap topology in both cases of synthetic and spatial dimensions. We start with the topology in one synthetic dimension. Our analysis reveals the reduction of $\mathbb{Z} \times \mathbb{Z} \to \mathbb{Z}$ for systems with charge U(1) symmetry and spin-parity symmetry. We end up this conclusion by analyzing a toy model, as well as by an argument in terms of topological invariants. Furthermore, we analyze an extended Hatano-Nelson chain where such reduction results in a striking phenomenon: fragility of a non-Hermitian skin effect against interactions in one spatial dimension.

The rest of this paper is organized as follows. In Sec. II, we discuss the reduction of non-Hermitian topological classification in one synthetic dimension by introducing topological invariants. In Sec. III, computing the obtained topological invariants, we address the reduction of topological classification in one spatial dimension. A brief summary and discussions are provided in Sec. IV. In Appendix A, detailed analysis of a non-Hermitian quantum dot is provided. Appendix B is devoted to detailed analysis of the extended Hatano-Nelson chain.

II. POINT-GAP TOPOLOGY IN ONE SYNTHETIC DIMENSION

A. Topological invariants

Firstly, we provide a generic argument in terms of topological invariants. Consider a quantum dot whose manybody Hamiltonian reads

$$\hat{H} = \hat{H}_0(\theta) + \hat{H}_{\text{int}},\tag{1}$$

with $\hat{H}_0(\theta) = \sum_{\alpha\beta} \hat{\Psi}^{\dagger}_{\alpha} h_{\alpha\beta}(\theta) \hat{\Psi}_{\beta}$, and $\hat{\Psi}^T = (\hat{c}_{a\uparrow}, \hat{c}_{a\downarrow}, \hat{c}_{b\uparrow}, \hat{c}_{b\downarrow}, \ldots)$. The second term \hat{H}_{int} denotes twobody interactions of fermions. Here, one-body Hamiltonian $h(\theta)$ is non-Hermitian and satisfies $h(2\pi) = h(0)$. The synthetic dimension is parameterized by θ which corresponds to a tunable parameter in experiments^{127,128} (e.g., a hopping integral in cold atoms). The operator $\hat{c}^{\dagger}_{l\sigma}$ ($\hat{c}_{l\sigma}$) creates (annihilates) a fermion in orbital l $(l = a, b, \ldots)$ and spin state σ ($\sigma = \uparrow, \downarrow$). The subscript α labels the set of l and σ .

Throughout this paper, we suppose that the Hamiltonian (1) respects the charge U(1) symmetry and spinparity symmetry. Namely, the zero-dimensional Hamiltonian satisfies

$$[\hat{H}, \hat{N}] = 0, \qquad [\hat{H}, e^{i\pi\hat{S}^z}] = 0,$$
 (2)

with $\hat{N} = \sum_{\alpha} \hat{\Psi}^{\dagger}_{\alpha} \hat{\Psi}_{\alpha}$ and $\hat{S}^z = \sum_{l=a,b,\dots} (\hat{c}^{\dagger}_{l\uparrow} \hat{c}_{l\uparrow} - \hat{c}^{\dagger}_{l\downarrow} \hat{c}_{l\downarrow})/2.$

Here, let us discuss the point-gap topology of the above system. In terms of the one-body Hamiltonian, we can introduce two distinct \mathbb{Z} -invariants due to the spin-parity symmetry. Because the one-body Hamiltonian $h(\theta)$ is periodic in θ , we can introduce the winding number w

$$w = \int_{0}^{2\pi} \frac{d\theta}{2\pi i} \partial_{\theta} \operatorname{tr} \log \left[h(\theta) - \epsilon_{\operatorname{ref}} \mathbb{1} \right], \qquad (3)$$

with the reference energy $\epsilon_{\text{ref}} \in \mathbb{C}$. The derivative with respect to θ is denoted by ∂_{θ} . The symbol "tr" denotes the trace of a matrix (i.e., $\operatorname{tr} h = \sum_{\alpha} h_{\alpha\alpha}$).

In addition, we can introduce spin winding number $w_{\rm s}$

$$w_{\rm s} = \int_0^{2\pi} \frac{d\theta}{4\pi i} \partial_\theta \mathrm{tr} \left[s^z \log \left(h(\theta) - \epsilon_{\rm ref} \mathbb{1} \right) \right], \qquad (4)$$

with $(s^z)_{\alpha\beta} = \operatorname{sgn}(\sigma)\delta_{\alpha\beta}$. Here, $\delta_{\alpha\beta}$ takes 1 (0) for $\alpha = \beta$ $(\alpha \neq \beta)$, and $\operatorname{sgn}(\sigma)$ takes 1 (-1) for $\sigma = \uparrow (\downarrow)$. For the spin winding number the spin-parity symmetry is essential; the one-body Hamiltonian satisfies $[s^z, h(\theta)] = 0$ in the presence of the spin-parity symmetry¹²⁹.

The above results indicates that in the presence of the U(1) symmetry and the spin-parity symmetry, the pointgap topology of $h(\theta)$ is characterized by two distinct \mathbb{Z} invariants.

Now, let us discuss the point-gap topology of the manybody Hamiltonian. In the presence of the spin-parity symmetry, the Hamiltonian \hat{H} can be block-diagonalized with \hat{N} and $\hat{P} := (-1)^{\hat{N}_{\uparrow}} = e^{i\frac{\pi}{2}\hat{N}}e^{i\pi\hat{S}^z}$. Here, \hat{N}_{\uparrow} denotes the operator of total number of fermions in the up-spin state. Therefore, for each Fock space, the following many-body winding number $W_{(N,P)}$ can be introduced¹³⁰;

$$W_{(N,P)}(E_{\mathrm{ref}}) = \int_0^{2\pi} \frac{d\theta}{2\pi i} \partial_\theta \mathrm{Tr} \log[\hat{H}_{(N,P)} - E_{\mathrm{ref}}\mathbb{1}], \ (5)$$

where N and P are eigenvalues of \hat{N} and \hat{P} , respectively. The reference energy is denoted by $E_{\text{ref}} \in \mathbb{C}$. By $\hat{H}_{(N,P)}$, we denote the many-body Hamiltonian for the subsector with (N, P). The symbol "Tr" denotes the trace over the subsector of the Fock space.

In the absence of interactions, eigenvalues of the manybody Hamiltonian $\hat{H}_{(N,P)}$ for each Fock space is computed from the eigenvalues of the one-body Hamiltonian $h(\theta)$ whose point-gap topology is characterized by w and $w_{\rm s}$.

The above results indicate that the point-gap topology of the one-body Hamiltonian $h(\theta)$ is characterized by a set of two Z-invariants (w, w_s) while the topology of the many-body Hamiltonian \hat{H} is characterized by the Zinvariant $W_{(N,P)}$ for each sector of the Fock space. As we see below, this fact is consistent with the behavior that the non-trivial topology characterized by $(w, w_s) =$ (0, 1) is trivialized by introducing the interactions.

B. Two orbital quantum dot: non-interacting case

As a specific case of Eq. (1), let us consider a twoorbital quantum dot (l = a, b) with a diagonal matrix $h(\theta) [h_{\alpha\beta}(\theta) = h_{\alpha}(\theta)\delta_{\alpha\beta}]$ whose diagonal elements are written as

$$h_{\alpha}(\theta) = \lambda e^{i\theta} \delta_{\alpha,(a,\uparrow)} + \lambda e^{-i\theta} \delta_{\alpha,(a,\downarrow)} + i\epsilon_{l\sigma} \delta_{\alpha,(l,\sigma)}.$$
 (6)

Here, λ and $\epsilon_{l\sigma}$ $(l = a, b \text{ and } \sigma = \uparrow, \downarrow)$ are real numbers. At the non-interacting level, couplings between orbitals are absent. The one-body Hamiltonian of orbital *a* corresponds to the small cycle limit of an extended Hatano-Nelson chain under the twisted boundary condition [see Eq. (10)].

The topology of $h(\theta)$ is characterized as $(w, w_s) = (0, 1)$ for $\epsilon_{ref} = 0$ and $|\epsilon_{a\sigma}| < \lambda$ ($\sigma = \uparrow, \downarrow$). To be concrete,



FIG. 1. Spectral flow of the one-body Hamiltonian $h(\theta)$. Data for the subsector with $\sigma = \uparrow [\sigma = \downarrow]$ are plotted in panel (a) [(b)]. The color denotes the value of θ . The data are obtained for $(\lambda, \epsilon_{a\uparrow}, \epsilon_{a\downarrow}, \epsilon_{b\uparrow}, \epsilon_{b\downarrow}) = (1, 0.2, -0.1, 0.35, -0.25).$

we plot a spectral flow of the one-body Hamiltonian in Fig. 1 for $(\lambda, \epsilon_{a\uparrow}, \epsilon_{a\downarrow}, \epsilon_{b\uparrow}, \epsilon_{b\downarrow}) = (1, 0.2, -0.1, 0.35, -0.25)$. This figure indicate that increasing θ from 0 to 2π , an eigenvalue winds around the origin in the clockwise (counter-clockwise) direction for the subsector $\sigma = \uparrow$ $(\sigma = \downarrow)$. The above numerical data support that the topology of $h(\theta)$ is characterized as $(w, w_s) = (0, 1)$. Figure 2(a) displays a spectral flow of the many-body Hamiltonian \hat{H}_0 for the subsector with (N, P) = (2, 1) of the Fock space. We can observe the loop structure of the spectral flow due to the topology of the one-body Hamiltonian $h(\theta)$. However, this figure indicates that $\hat{H}_{(2,1)}$ is topologically trivial [i.e., $W_{(2,1)} = 0$] for $E_{\text{ref}} = 0$ because an eigenvalue winds around the origin in the clockwise direction, and the other eigenvalue winds around the origin in the opposite direction.

C. Two orbital quantum dot: interacting case

Now, let us introduce the following two-body interaction

$$\hat{H}_{\rm int} = \frac{iJ}{2} (\hat{S}_a^+ \hat{S}_b^- + \text{h.c.}) + \frac{iV}{2} (\hat{S}_a^+ \hat{S}_b^+ + \text{h.c.}), \quad (7)$$

with real numbers J and V. Here, "h.c." denotes the Hermitian conjugate of the corresponding operator [e.g., $iJ(\hat{S}_a^+\hat{S}_b^- + h.c.) = iJ(\hat{S}_a^+\hat{S}_b^- + \hat{S}_a^-\hat{S}_b^+)$]. The spin operator \hat{S}_l^\pm is defined as $\hat{S}_l^\pm = \hat{S}_l^x \pm i\hat{S}_l^y$ with $\hat{S}_l^{x(y)}$ being the x- (y-) component of the spin operator for orbital l. The above two-body interactions respect charge U(1) symmetry and spin-parity symmetry; applying the operator $e^{i\pi\hat{S}^z}$ transforms the spin operators as $e^{i\pi\hat{S}^z}\hat{S}_l^\pm e^{-i\pi\hat{S}^z} = -\hat{S}_l^\pm$, meaning that the interactions respect spin-parity symmetry.

For the sake of simplicity, we focus on the subsector with (N, P) = (2, 1). The results for the subsector (N, P) = (2, -1) are provided in Appendix A 1. Figure 2(b) displays the spectral flow for V = J = 1. Remarkably, this figure indicate that the interactions open a imaginary gap; interactions split the loops which wind the origin at the non-interacting level [see Fig. 2(a)].

This fact indicates that interactions [Eq. (7)] allow a smooth deformation of the spectral flow for $\lambda = 1$ to that for $\lambda = 0$ without closing the point-gap at $E_{\text{ref}} = 0$ the latter of which is obviously trivial. Indeed, the following deformation smoothly connects the Hamiltonian $\hat{H}(\theta)$ for $\lambda = 1$ and that for $\lambda = 0$: (i) Increasing V from 0 to 1 for $\lambda = 1$ and J = V [see Fig. 2(c)]; (ii) Decreasing λ from 1 to 0 for $J = V = \sqrt{\lambda}$ [see Fig. 2(d)]. This deformation demonstrates that the many-body Hamiltonian $\hat{H}_{(2,1)}(\theta)$ is topologically trivial.

We note that difference of the symmetry constraint of the spin-parity symmetry is essential for the imaginary gap at ImE = 0 in Fig. 2(b). As discussed above, the symmetry constraints (2), which results in $[s^z, h(\theta)] = 0$, forbids hybridization terms between two distinct subsectors with (N, S^z) . In contrast, the symmetry constraint allows such hybridization terms of two-body interactions \hat{H}_{int} . Therefore, the two-body interactions can destroy the loop structure arising from the non-trivial topology of the one-body Hamiltonian [see Figs. 2(a) and 2(b)].

For instance, in the subsector with (N, P) = (2, 1), the



FIG. 2. Spectral flow of the many-body Hamiltonian for the subsector with (N, P) = (2, 1). (a): Spectral flow for J = V = 0 and $\lambda = 1$. Data for the subsector with $(N, S^z) = (2, 1)$ [(2, -1)] are plotted with open circles [closed triangles]. (b): Spectral flow for $V = J = \lambda = 1$. In panels (a) and (b), we can see that the eigenvalues flow as denoted by arrows with increasing θ from 0 to 2π . (c): Spectral flow for several values of V (J = V) at $\lambda = 1$. With increasing V from 0 to 1, the eigenvalues flow as denoted by arrows. (d): Spectral flow for several values of λ for $V = J = \sqrt{\lambda}$. With decreasing λ from 1 to 0, the eigenvalues flow as denoted by arrows. The data are obtained for $(\epsilon_{a\uparrow}, \epsilon_{a\downarrow}, \epsilon_{b\uparrow}, \epsilon_{b\downarrow}) = (0.2, -0.1, 0.35, -0.25)$.

Hamiltonian is written as

$$\hat{H}_{(2,1)} = \begin{pmatrix} \lambda e^{i\theta} + i\epsilon_{a\uparrow} + i\epsilon_{b\uparrow} & \frac{iV}{2} \\ \frac{iV}{2} & \lambda e^{-i\theta} + i\epsilon_{a\downarrow} + i\epsilon_{b\downarrow} \end{pmatrix}.$$

Here, we have chosen the following basis vectors spanning the subsector of the Fock space $(\hat{c}^{\dagger}_{a\uparrow}\hat{c}^{\dagger}_{b\uparrow}|0\rangle, \hat{c}^{\dagger}_{a\downarrow}\hat{c}^{\dagger}_{b\downarrow}|0\rangle)$. The vacuum state is denoted by $|0\rangle$ (i.e., $\hat{c}_{l\sigma}|0\rangle = 0$ for arbitrary l and σ).

Diagonalizing the above Hamiltonian, we obtain

$$E_{\pm} = \lambda \cos \theta + i\delta_0 \pm i\sqrt{(\sin \theta + \delta_3)^2 + (\frac{V}{2})^2}, \quad (8)$$

with $2\delta_0 = \epsilon_{a\uparrow} + \epsilon_{b\uparrow} + \epsilon_{a\downarrow} + \epsilon_{b\downarrow}$ and $2\delta_3 = \epsilon_{a\uparrow} + \epsilon_{b\uparrow} - \epsilon_{a\downarrow} - \epsilon_{b\downarrow}$. Equation (8) elucidates that spin-parity symmetry allows the hybridization term between states with $(N, S^z) = (2, 1)$ and $(N, S^z) = (2, -1)$ which opens the line-gap $\text{Im}[E_+(\theta) - E_-(\theta)] > 0$ [see Fig. 2(b)]. In contrast, spin-parity symmetry forbids such hybridization terms for the quadratic Hamiltonian \hat{H}_0 .

The above numerical results supports that the manybody Hamiltonian $\hat{H}_{(2,1)}(\theta)$ is topologically trivial despite the loop structure due to the topology of the onebody Hamiltonian with $(w, w_s) = (0, 1)$. We can also confirm the robustness of the topology characterized by finite values of $W_{(2,1)}^{131}$ (see also Appendix A 2). Putting the argument in terms of the topological invariants and the above results of the toy model together, we end up with the reduction of the point-gap topology $\mathbb{Z} \times \mathbb{Z} \to \mathbb{Z}$.

III. POINT-GAP TOPOLOGY IN ONE SPATIAL DIMENSION AND FRAGILITY OF A NON-HERMITIAN SKIN EFFECT

By analyzing an extended Hatano-Nelson chain [see Fig. 3(a)], we elucidate that interactions reduce the point-gap topology in one spatial dimension as is the case in one synthetic dimension. Remarkably, this reduction phenomenon results in fragility of a non-Hermitian skin effect against interactions. As in the case of one synthetic dimension, essential ingredients are spin-parity symmetry and two-body terms flipping spins.

Let us consider an extended Hatano-Nelson chain [see Fig. 3(a)] whose Hamiltonian reads

$$\hat{H}_{\rm eHN}(\theta) = \hat{H}_0(\theta) + \hat{H}_{\rm int}, \qquad (9a)$$

$$\hat{H}_0(\theta) = \sum_k \hat{\Psi}_{k\alpha}^{\dagger} h_{\alpha\beta}(k,\theta) \hat{\Psi}_{k\beta}, \qquad (9b)$$

$$\hat{H}_{\text{int}} = \sum_{j=0,L-1} \left[J(\hat{S}_{ja}^{+} \hat{S}_{jb}^{-} + \text{h.c.}) + iV(\hat{S}_{ja}^{+} \hat{S}_{jb}^{+} + \text{h.c.}) \right],$$
(9c)

with a diagonal matrix $h(k,\theta)$ $[h_{\alpha\beta}(k,\theta) = h_{\alpha}(k,\theta)\delta_{\alpha\beta}, (kL/2\pi = 0, 1, ..., L-1)]$ whose diagonal elements are

$$h_{\alpha}(k,\theta) = t\delta_{\alpha,(a,\uparrow)}e^{i(k+\theta/L)} + t\delta_{\alpha,(a,\downarrow)}e^{-i(k+\theta/L)}.(10)$$

Here, we have imposed the twisted boundary condition in order to compute the winding numbers (for more details, see Appendix B 1). The operator $\hat{\Psi}_{k\alpha}$ is the Fourier transformed annihilation operator $\Psi_{k\alpha} := \frac{1}{\sqrt{L}} \sum_{j=0,\dots,L-1} e^{ikj} \Psi_{j\alpha}$ with $\Psi_j^T = (\hat{c}_{ja\uparrow}, \hat{c}_{ja\downarrow}, \hat{c}_{jb\uparrow}, \hat{c}_{jb\downarrow})$. The two-body term \hat{H}_{int} describes the interaction between fermions in orbital *a* and localized fermions in orbital *b*. This model also preserves charge U(1) and spinparity symmetry, meaning that $\hat{H}_{eHN}(\theta)$ can be block-diagonalized with \hat{N} and $\hat{P} = (-1)^{\hat{N}_{a\uparrow} + \hat{N}_{b\uparrow}}$ where $\hat{N}_{l\sigma}$ and \hat{N} are defined as $\hat{N}_{l\sigma} = \sum_{j} \hat{c}_{jl\sigma}^{\dagger} \hat{c}_{jl\sigma}$ and $\hat{N} = \sum_{l\sigma} \hat{N}_{l\sigma}$, respectively. The Hamiltonian $\hat{H}_{eHN}(\theta)$ also commutes with $\hat{n}_{jb} = \sum_{\sigma} \hat{c}_{jb\sigma}^{\dagger} \hat{c}_{jb\sigma}$ for j = 0, L - 1, and thus, we suppose that orbital *b* is occupied at both edges (j = 0, L - 1) by focusing on the corresponding Fock space.

Now, we demonstrate that for $N_a = 1$ (i.e., N = 3), a non-Hermitian skin effect observed at the non-interacting level is fragile against the two-body interactions due to trivial topology of the many-body Hamiltonian. Let us start with the non-interacting level. Under the twisted boundary condition, the spectral flow shows a loop structure [see Fig. 3(b)] due to the point-gap topology of onebody Hamiltonian $h(\theta) := \bigoplus_k h(k, \theta)$ characterized by



FIG. 3. (a): Sketch of the extended Hatano-Nelson chain. (b) [(d)]: Spectral flow for J = V = 0 [J = V = 1]. Red dots denote the data obtained under the open boundary condition. (c) [(e)]: Expectation values of $\hat{n}_{ja\sigma}$ for J = V = 0[J = V = 1] under the open boundary condition. In panels (c) and (e), $\langle \hat{n}_{ja\sigma} \rangle = {}_{\rm R} \langle \Phi_n | \hat{n}_{ja\sigma} | \Phi_n \rangle_{\rm R}$ is plotted against j and $\langle \hat{N}_{a\uparrow} \rangle = {}_{\rm R} \langle \Phi_n | \hat{N}_{a\uparrow} | \Phi_n \rangle_{\rm R}$. Here, $|\Phi_n \rangle_{\rm R}$ (n = 0, 1, ...) denote right eigenstates of $\hat{H}_{\rm eHN}$. Red (blue) lines denote the data for $\sigma = \uparrow (\sigma = \downarrow)$. These data are obtained for the subsector (N, P) = (3, -1) and a parameter set (L, t) = (7, 1).

 $(w, w_s) = (0, 1)$ for $\epsilon_{ref} = 0$. This non-trivial topology of *h* induces the non-Hermitian skin effect at the noninteracting level. In the presence of the boundaries, all of the eigenvalues $[E_n \ (n = 0, 1, ...)]$ become zero in contrast to the eigenvalues in the absence of the boundaries [see Fig. 3(b)]. In addition, a fermion in the up- (down-) spin state is localized around the right (left) edge under the open boundary condition [see Fig. 3(c)].



FIG. 4. Expectation values of $\hat{n}_{ja\uparrow}$ for the subsector (N, P) = (9, -1) and a parameter set (L, t) = (7, 1). In panels (a) and (b), data for J = V = 0 and J = V = 1 are plotted, respectively. These figures are plotted in the same way as Figs. 3(c) and 3(e).

However, interactions destroy the above non-Hermitian skin effect, which is due to the trivial topology of the many-body Hamiltonian $W_{(3,-1)} = 0$ for $E_{\rm ref} = 0$ (for computation of the many-body winding number, see Fig. 6). Because of the trivial topology, we can observe that interactions destroy the loop structure of the spectral flow and open a line-gap for J = V = 1[see Fig. 3(d)], which is also confirmed by analysis based on the perturbation theory (see Appendix B2). This result verifies the reduction of the point-gap topology $\mathbb{Z} \times \mathbb{Z} \to \mathbb{Z}$ for the subsector with (N, P) = (3, -1). Correspondingly, the interactions destroy the extreme sensitivity of the spectrum to the presence/absence of boundaries [see Fig. 3(d)]. Furthermore, in the presence of interactions fermions extend to the bulk even under the open boundary condition [see Fig. 3(e)]. This result is also intuitively understood as follows: while the one-body term $\hat{H}_0(\theta)$ localizes the fermions in orbital a and the up- (down-) spin state around the right (left) edge, the two-body interactions \hat{H}_{int} flip their spins at edges, which suppresses the effects of boundaries. The above results indicate that the non-Hermitian skin effect observed at the non-interacting level is fragile against the two-body interactions¹³². Our numerical calculations indicate that such fragility of the non-Hermitian skin effect is also observed for the case of many fermions in orbital a [see Fig. 4]. More detailed data are provided in Appendix B 3.

IV. SUMMARY AND DISCUSSION

We have analyzed interaction effects on the onedimensional point-gap topology in both cases of synthetic and spatial dimensions. Our analysis has elucidated that the reduction $\mathbb{Z} \times \mathbb{Z} \to \mathbb{Z}$ occurs for systems of synthetic one dimension with charge U(1) symmetry and spin-parity symmetry. This conclusion is obtained by the argument of topological invariants as well as by explicit analysis of the toy model. Furthermore, we have also analyzed the extended Hatano-Nelson chain which exhibits striking interaction effects: interactions reduce the point-gap topology and destroy the non-Hermitian skin effect at the non-interacting level.

We stress that the spin-parity symmetry plays an essential role, which leads qualitative differences from results of previous works^{119,120}. Instead of Eq. (9c), one can introduce the Hubbard type interactions preserving spin U(1) symmetry. This type of interactions does not flip the spins and thus would not destroy the non-Hermitian skin effect as discussed in Refs. 119 and 120. Analysis for other types of interactions is left as future works to be addressed. We also remark several open questions. Our results indicate that interactions result in the same reduction phenomenon for both cases of synthetic and spatial dimension whose generality for other symmetry classes remains as an open question. In addition, establishing strict one to one correspondence between the many-body winding number and skin effect also remains a crucial open question.

The above discoveries shed new light on non-Hermitian correlated systems and open up a new directions of researches on non-Hermitian topological physics. For instance, the above results imply the possibility of similar reduction phenomena for other cases of symmetry and dimensions. As well as the above theoretical open question, experimental observation of the reduction is also a significant issue to be addressed. We expect that cold atoms are promising candidate where interactions and non-Hermiticity can be tuned in experiments.

V. ACKNOWLEDGMENTS

This work was supported by MEXT KAKENHI Grantin-Aid for Transformative Research Areas A "Extreme Universe" No. JP22H05247. This work is also supported by JSPS KAKENHI Grant No. JP21K13850 and also by JST CREST, Grant No. JPMJCR19T1, Japan.

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- ¹²⁷ In the literatures, there are two types of synthetic dimensions. Reference 35 regards internal degrees of freedoms of particles as synthetic dimension, while Ref. 33 regards periodic parameters as synthetic dimensions. By making use of the latter type of synthetic dimensions, twodimensional topology has been studied for a system in one spatial dimension^{36,37}. Following these works, we regard the parameter θ as a synthetic dimension and discuss onedimensional non-Hermitian topology for the quantum dot.
- ¹²⁸ In the case of synthetic dimension, one does not have to consider thermodynamic limit which significantly simplifies the analysis.
- ¹²⁹ Instead of w and $w_{\rm s}$, one can also use w_{\uparrow} and w_{\downarrow} which are rewritten as $2w_{\uparrow} = w + 2w_{\rm s}$ and $2w_{\downarrow} = w - 2w_{\rm s}$. The winding number $w_{\uparrow} (w_{\downarrow})$ can be computed in a similar way as Eq. (3) by replacing h to the one-body Hamiltonian for the up-spin (down-spin) sector $h_{\uparrow} (h_{\downarrow})$.
- ¹³⁰ We note that introducing additional terms makes other subsectors irrelevant. For instance adding the term μ(N̂ − 2) with μ ∈ C to Eq. (1) can make subsectors labeled by N = 0, 1, 3, 4 irrelevant to the topology for E_{ref} = 0.
- ¹³¹ Equation (5) indicates that the topology characterized by a finite value of the many-body winding number is robust against interactions. Its robustness can be explicitly checked for $W_{(2,1)} = 1$ and $W_{(2,1)} = 2$. For $W_{(2,1)} = 1$, the robustness of the topology can be seen in Figs. 2(a) and 2(b). Figure 2(a) indicates that the topology is characterized by $W_{(2,1)} = 1$ for $E_{\text{ref}} = i$ in the absence of the interactions. Figure 2(b) indicates that this nontrivial topology is maintained in the presence of the interactions. The topology characterized by $W_{(2,1)} = 2$ can be discussed by replacing h whose robustness is discussed in Appendix A 2.
- ¹³² At the non-interacting level, fragility of the skin effect against spin-flipping terms has been discussed in Ref. 55. In the previous work, the one-body term maintains the point-gap topology nontrivial, while our two-body interactions destroy the nontrivial topology.

Appendix A: Details of a two-orbital model

1. Analysis for the subsector with (N, P) = (2, -1)

In the main text, we have seen that interactions open a line-gap for the subsector with (N, P) = (2, 1), which is consistent with the trivial topology $W_{(2,1)} = 0$ for $E_{\text{ref}} =$ 0. In this section, we show that a similar behavior is observed for the subsector with (N, P) = (2, -1).



FIG. 5. Spectral flow of the many-body Hamiltonian for the subsector with (N, P) = (2, -1). Panels (a)-(d) are plotted in the same way as panels (a)-(d) in Fig. 2. Panel (e) is a magnified version of the range $-0.12 \leq \text{Im}E \leq 0.12$ in panel (d).

Despite the non-trivial topology of the one-body Hamiltonian, the many-body winding number takes zero $[W_{(2,-1)} = 0]$ for $E_{\text{ref}} = 0$ as shown in Fig. 5(a). Correspondingly, the spectrum of the many-body Hamiltonian $\hat{H}_{(2,-1)}$ can smoothly shrink to the points [see Figs. 5(b)-5(d)].

In this subsector, the interaction J is essential for destruction of the loop structure observed in Fig. 5(a), which can be seen as follows. In the subsector with (N, P) = (2, -1), the Hamiltonian is written as

$$H_{(2,-1)} = H_{0(2,-1)} + H_{int(2,-1)},$$
 (A1a)

$$\hat{H}_{0(2,-1)} = \operatorname{diag} \left(\lambda e^{i\theta} + i\epsilon_{a\uparrow} + i\epsilon_{b\downarrow}, \lambda e^{-i\theta} + i\epsilon_{a\downarrow} + i\epsilon_{b\uparrow} \\ 2\lambda \cos\theta + i\epsilon_{a\uparrow} + i\epsilon_{a\downarrow}, i\epsilon_{b\uparrow} + i\epsilon_{b\downarrow} \right), \quad (A1b)$$

with $diag(\cdots)$ denoting a diagonal matrix. Here, we have chosen the following basis vectors spanning the subsector of the Fock space

$$\left(\hat{c}^{\dagger}_{a\uparrow}\hat{c}^{\dagger}_{b\downarrow}|0\rangle,\hat{c}^{\dagger}_{a\downarrow}\hat{c}^{\dagger}_{b\uparrow}|0\rangle,\hat{c}^{\dagger}_{a\uparrow}\hat{c}^{\dagger}_{a\downarrow}|0\rangle,\hat{c}^{\dagger}_{b\uparrow}\hat{c}^{\dagger}_{b\downarrow}|0\rangle\right).$$
 (A2)

Diagonalizing the Hamiltonian, we obtain

$$E_{\pm} = \lambda \cos \theta + i\delta_0' \pm i\sqrt{(\sin \theta + \delta_3')^2 + (\frac{J}{2})^2}, \quad (A3)$$

$$E' = 2\lambda\cos\theta + i\epsilon_{a\uparrow} + i\epsilon_{a\downarrow},\tag{A4}$$

$$E'' = i\epsilon_{b\uparrow} + i\epsilon_{b\downarrow},\tag{A5}$$

with $2\delta'_0 = \epsilon_{a\uparrow} + \epsilon_{b\downarrow} + \epsilon_{a\downarrow} + \epsilon_{b\uparrow}$ and $2\delta'_3 = \epsilon_{a\uparrow} + \epsilon_{b\downarrow} - (\epsilon_{a\downarrow} + \epsilon_{b\uparrow})$. The above results elucidate that the spinparity symmetry allows the two-body interaction which splits loop structure observed in Fig. 5(a).

2. Robustness of the topology characterized by $W_{(2,1)}=2 \label{eq:W2}$

Let us analyze a toy model in order to demonstrate the robustness of the topology characterized by $W_{(2,1)} = 2$. Specifically, consider \hat{H} specified by

$$h_{\alpha}(\theta) = \lambda e^{i\theta} (\delta_{\alpha,(a,\uparrow)} + \delta_{\alpha,(a,\downarrow)}) + i\epsilon_{l\sigma} \delta_{\alpha,(l,\sigma)},$$
 (A6)

and the interaction term [Eq. (7)]. Then, we have

$$\hat{H}_{(2,1)} = \begin{pmatrix} \lambda e^{i\theta} + i\epsilon_{a\uparrow} + i\epsilon_{b\uparrow} & \frac{iV}{2} \\ \frac{iV}{2} & \lambda e^{i\theta} + i\epsilon_{a\downarrow} + i\epsilon_{b\downarrow} \end{pmatrix},$$

for the subsector with (N, P) = (2, 1) [see also arguments above Eq. (8)]. Diagonalizing the above Hamiltonian, we obtain

$$E_{\pm} = \lambda e^{i\theta} + i\delta_0 \pm i\sqrt{(\delta_3)^2 + (\frac{V}{2})^2},$$
 (A7)

where δ_0 and δ_3 are defined just below Eq. (8). This result demonstrate that the topology characterized by $W_{(2,1)} = 2$ is robust against interactions. Let us choose the parameters as $\delta_0 = \delta_3 = 0$. In this case, Eqs. (5) and (A7) indicate that the topology is characterized by $W_{(2,1)} = 2$ for $E_{\text{ref}} = 0$ in the absence of the interactions. Equation (A7) also indicates that this nontrivial topology is maintained for finite values of V.

Appendix B: Details of the extended Hatano-Nelson chain

1. Hamiltonian under the twisted boundary condition

We provide the explicit form of the extended Hatano-Nelson chain under the twisted boundary condition. The Hamiltonian reads

$$\hat{H}_{\rm eHN} = \hat{H}_0(\theta) + \hat{H}_{\rm int},\tag{B1a}$$

$$\hat{H}_{0}(\theta) = t \left[e^{i\theta} \hat{c}^{\dagger}_{0a\uparrow} \hat{c}_{L-1a\uparrow} + \sum_{j=0}^{L-2} \hat{c}^{\dagger}_{j+1a\uparrow} \hat{c}_{ja\uparrow} \right] + t \left[e^{-i\theta} \hat{c}^{\dagger}_{L-1a\downarrow} \hat{c}_{0a\downarrow} + \sum_{j=1}^{L-1} \hat{c}^{\dagger}_{j-1a\downarrow} \hat{c}_{ja\downarrow} \right],$$
(B1b)

$$\hat{H}_{\rm int}(\theta) = \sum_{j=0,L-1} \left[\frac{J}{2} (\hat{S}_{ja}^+ \hat{S}_{jb}^- + \hat{S}_{ja}^- \hat{S}_{jb}^+) + iV (\hat{S}_{ja}^+ \hat{S}_{jb}^+ + \hat{S}_{ja}^- \hat{S}_{jb}^-) \right].$$
(B1c)

Under a gauge transformation $\hat{c}_{ja\sigma} \to e^{-i\frac{\theta}{L}j}\hat{c}_{ja\sigma}$, the one-body term is written as

$$\hat{H}_{0}(\theta) = \sum_{j=0}^{L-1} \left[t e^{i\theta/L} \hat{c}^{\dagger}_{j+1a\uparrow} \hat{c}_{ja\uparrow} + t e^{-i\theta/L} \hat{c}^{\dagger}_{j-1a\downarrow} \hat{c}_{ja\downarrow} \right],$$
(B2)

with $\hat{c}^{\dagger}_{La\uparrow} := \hat{c}^{\dagger}_{0a\uparrow}$ and $\hat{c}^{\dagger}_{-1a\downarrow} := \hat{c}^{\dagger}_{L-1a\downarrow}$.

Applying the Fourier transformation to the above Hamiltonian yields Eq. (10). We note that under the

open boundary condition, hopping terms between sites j = 0 and j = L - 1 [i.e., the first and the third terms of Eq. (B1b)] become zero.

In the presence of charge U(1) symmetry and spinparity symmetry, the one-body Hamiltonian is characterized by w and w_s [Eqs. (3) and (4)] with $h(\theta) := \bigoplus_k h(k, \theta)$. The topology of the many-body Hamiltonian for given subsector with (N, P) is characterized by the many-body winding number [Eq. (5)] with $\hat{H}_{(N,P)} :=$ $\hat{H}_{eHN(N,P)}$. Here, $\hat{H}_{eHN(N,P)}$ denotes the many-body Hamiltonian of the extended Hatano-Nelson model for the given subsector with (N, P).

2. Analysis based on the perturbation theory

Based on the perturbation theory, we confirm that interactions open a line-gap as shown in Fig. 3(d). As mentioned in the main text, we suppose that orbital b is occupied at both edges (j = 0, L - 1).

Suppose that interactions are sufficiently weak. In the subsector of (N, P) = (3, -1), the non-interacting Hamiltonian $\hat{H}_0(\theta)$ is written as

$$\hat{H}_{0(3,-1)} = t\omega^n \begin{pmatrix} e^{i\frac{\theta}{L}} & 0 & 0 & 0\\ 0 & e^{i\frac{\theta}{L}} & 0 & 0\\ 0 & 0 & e^{-i\frac{\theta}{L}} & 0\\ 0 & 0 & 0 & e^{-i\frac{\theta}{L}} \end{pmatrix},$$
(B3)

with $\omega = e^{\frac{2\pi i}{L}}$ and the basis

$$(|n\uparrow;\uparrow\uparrow\rangle,|n\uparrow;\downarrow\downarrow\rangle,|n\downarrow;\uparrow\downarrow\rangle,|n\downarrow;\downarrow\uparrow\rangle), \tag{B4}$$

for given n (n = 0, 1, 2, ..., L - 1). Here, $|n\sigma; \sigma'\sigma''\rangle$ is defined as $|n\sigma; \sigma'\sigma''\rangle := \overline{\hat{d}}_{n\sigma} \hat{c}^{\dagger}_{0b\sigma'} \hat{c}^{\dagger}_{L-1b\sigma''} |0\rangle$ with $\overline{\hat{d}}_{n\uparrow} := \sum_{j} \hat{c}^{\dagger}_{ja\uparrow} R_{jn}$ and $\overline{\hat{d}}_{n\downarrow} := \sum_{j} \hat{c}^{\dagger}_{ja\downarrow} L_{jn}^*$. Matrices R and L^{\dagger} $(R_{jn} := \frac{1}{\sqrt{L}} \omega^{-nj})$ and $L_{nj}^{\dagger} := \frac{1}{\sqrt{L}} \omega^{nj}$ diagonalize the matrix h $(h_{ij} = t\delta_{i,j+1})$

$$L^{\dagger}hR = t\operatorname{diag}(1,\omega,\omega^2,\cdots,\omega^{L-1})$$
(B5)

which corresponds to the kinetic term of fermions in orbital a and the up-spin state for $\theta = 0$ [see Eq. (B1b) and Fig. 3(a)]. Here diag(\cdots) describes a diagonal matrix. Introducing the operators $\hat{d}_{n\uparrow} := \sum_{j} \hat{c}_{ja\uparrow}(L^{\dagger})_{nj}$ and $\hat{d}_{n\downarrow} := \sum_{j} \hat{c}_{ja\downarrow}(R^T)_{nj}$, we have anti-commutation relations

$$\{\hat{d}_{n\sigma}, \hat{d}_{m\sigma'}\} = \delta_{nm} \delta_{\sigma\sigma'},\tag{B6}$$

for $n, m = 0, 1, 2, \dots, L - 1$ and $\sigma, \sigma' = \uparrow, \downarrow$, which can be seen by noting the relations $\sum_j L_{nj}^{\dagger} R_{jm} = \delta_{nm}$ and $\{\hat{c}_{il\sigma}, \hat{c}_{jl'\sigma'}^{\dagger}\} = \delta_{ij}\delta_{ll'}\delta_{\sigma\sigma'}$. We note that $(\bar{d}_{n\sigma})^{\dagger} = d_{n\sigma}$ holds due to the relation $R_{jn}^* = L_{nj}^{\dagger}$.

Now, let us compute energy eigenvalues at the first order of the interactions. Firstly, we note the following relations.

$$\hat{S}_{ja}^{+}|m\downarrow;\sigma\sigma'\rangle = \sum_{i} \hat{S}_{ja}^{+} L_{im}^{*} \hat{c}_{ia\downarrow}^{\dagger} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= L_{jm}^{*} \hat{c}_{ja\uparrow}^{\dagger} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= \sum_{n} L_{jm}^{*} (L^{\dagger})_{nj} \overline{\hat{d}}_{n\uparrow} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= \frac{1}{L} \sum_{n} \omega^{(n+m)j} |n\uparrow;\sigma\sigma'\rangle,$$
(B7)
$$\hat{S}_{ja}^{-}|m\uparrow;\sigma\sigma'\rangle = \sum_{i} \hat{S}_{ja}^{-} R_{im} \hat{c}_{ia\uparrow}^{\dagger} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= R_{jm} \hat{c}_{ja\downarrow}^{\dagger} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= \sum_{n} R_{jm} (R^{T})_{nj} \overline{\hat{d}}_{n\downarrow} \hat{c}_{0b\sigma}^{\dagger} \hat{c}_{L-1b\sigma'}^{\dagger}|0\rangle
= \frac{1}{L} \sum_{n} \omega^{-(n+m)j} |n\downarrow;\sigma\sigma'\rangle.$$
(B8)

Here, we have used the relations $\sum_{n} R_{jn} L_{ni}^{\dagger} = \delta_{ij}$, $\hat{c}_{ia\uparrow}^{\dagger} = \sum_{n} L_{ni}^{\dagger} \bar{\hat{d}}_{n\uparrow}$, and $\hat{c}_{ia\downarrow}^{\dagger} = \sum_{n} R_{ni}^{T} \bar{\hat{d}}_{n\downarrow}$. Thus, at the first order, the Hamiltonian is written as $\hat{H}_{eHN(3,1)} = \hat{H}_{0(3,1)} + \hat{H}_{int(3,1)}$ with $\hat{H}_{0(3,1)}$ in Eq. (B3) and

$$\hat{H}_{\text{int}(3,-1)} = \frac{V}{L} \begin{pmatrix} 0 & 0 & \omega^{-2n} & 1\\ 0 & 0 & 0 & 0\\ \omega^{2n} & 0 & 0 & 0\\ 1 & 0 & 0 & 0 \end{pmatrix} + \frac{iJ}{L} \begin{pmatrix} 0 & 0 & 0 & 0\\ 0 & 0 & 1 & \omega^{-2n}\\ 0 & 1 & 0 & 0\\ 0 & \omega^{2n} & 0 & 0 \end{pmatrix},$$
(B9)

for the basis defined in Eq. (B4).

The eigenvalues of $\hat{H}_{eHN(3,1)}$ are written as

$$E_{\rm p,\pm} = t\omega^n \left(\cos(\frac{\theta}{L}) \pm \sqrt{C_{\rm p}^2 - \sin^2(\frac{\theta}{L})} \right), \tag{B10a}$$

$$E_{\mathrm{m,\pm}} = t\omega^n \left(\cos(\frac{\theta}{L}) \pm \sqrt{C_{\mathrm{m}}^2 - \sin^2(\frac{\theta}{L})} \right), \tag{B10b}$$

with

$$C_{\rm p}^2 = \frac{1}{(t\omega^n L)^2} \left[(V^2 - J^2) + \sqrt{V^4 + J^4 - 2V^2 J^2 \text{Re}[\omega^{4n}]} \right],$$
(B10c)

$$C_{\rm m}^2 = \frac{1}{(t\omega^n L)^2} \left[(V^2 - J^2) - \sqrt{V^4 + J^4 - 2V^2 J^2 \text{Re}[\omega^{4n}]} \right].$$
 (B10d)

(B10e)

As well as by directly diagonalizing the matrix, the eigenvalues are obtained by taking square of the matrices (see below). These results indicate that interactions lift four-fold degeneracy observed for $\theta = 0$. Specifically, the imaginary parts of C_p and C_m lift the degeneracy. To see this, firstly, let us suppose that the imaginary parts are zero (Im $C_p = ImC_m = 0$), then, Eq. (B10a) indicates that exceptional points emerge at certain θ [i.e., $C_p^2 \ge 0$ holds, and $C_p^2 - ImC_m = 0$). $\sin^2(\theta/L) = 0$ can be satisfied]. On the other hand, the finite imaginary parts lift the degeneracy at $\theta = 0$ without inducing exceptional points [i.e., $C_p^2 - \sin^2(\theta/L) \neq 0$ for $0 \le \theta < 2\pi$]. Equations (B10c) and (B10d) indicate that the imaginary parts of C_p and C_m can be finite for proper choice of n,

J, and V. Therefore, the above result of the first-order perturbation theory indicate that interactions open a line-gap. We show that eigenvalues (B10) can be obtained by taking squares of the matrices. Consider the following matrix

$$\dot{H} = (x\sigma_0\tau_0 + y\sigma_0\tau_3 + a\sigma_1\tau_1 + b\sigma_2\tau_2 + c\sigma_0\tau_1 + d\sigma_3\tau_1 + f\sigma_0\tau_2 + g\sigma_3\tau_2),$$
(B11)

with complex numbers x, y, a, b, c, d, f and g. Here σ_0 and τ_0 denote the 2 × 2-identity matrix. Pauli matrices are denoted by σ_s and τ_s (s = 1, 2, 3). Matrices $\sigma_{\mu}\tau_{\nu}$ ($\mu, \nu = 0, 1, 2, 3$) denote 4×4 -matrices. For instance, $\sigma_1\tau_2$ is written as

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$$\sigma_1 \tau_2 = \begin{pmatrix} 0 & 0 & 0 & -i \\ 0 & 0 & -i & 0 \\ 0 & i & 0 & 0 \\ i & 0 & 0 & 0 \end{pmatrix}.$$
 (B12)

For the following parameter set,

$$\begin{pmatrix} x \\ y \\ a \\ b \\ c \\ d \\ f \\ g \end{pmatrix} = \frac{1}{2L} \begin{pmatrix} 2Lt\omega^{n}\cos\frac{\theta}{L} \\ 2iLt\omega^{n}\sin\frac{\theta}{L} \\ V+iJ \\ -V+iJ \\ (V+iJ)\text{Re}(\omega^{2n}) \\ (V-iJ)\text{Re}(\omega^{2n}) \\ (V-iJ)\text{Re}(\omega^{2n}) \\ (V-iJ)\text{Im}(\omega^{2n}) \end{pmatrix},$$
(B13)

 \tilde{H} is reduced to the matrix $\hat{H}_{eHN(3,-1)}$.

Taking square of this matrix yields

$$(\tilde{H} - x\sigma_0\tau_0)^2 - (y^2 + a^2 + b^2 + c^2 + d^2 + f^2 + g^2)$$

= $-2ab\sigma_3\tau_3 + 2(cd + fg)\sigma_3\tau_0 + 2ac\sigma_1\tau_0 + 2ag\sigma_2\tau_3 + 2bd\sigma_1\tau_3 + 2bf\sigma_2\tau_0.$ (B14)

Thus, we have

$$\begin{bmatrix} (\tilde{H} - x\sigma_0\tau_0)^2 - (y^2 + a^2 + b^2 + c^2 + d^2 + f^2 + g^2)^2 \end{bmatrix}^2$$

$$= \left[\{ -2ab\sigma_3\tau_3 + 2(cd + fg)\sigma_3\tau_0 \} + (2ac\sigma_1\tau_0 + 2bd\sigma_1\tau_3) + (2ag\sigma_2\tau_3 + 2bf\sigma_2\tau_0) \right]^2$$

$$= \{ -2ab\sigma_3\tau_3 + 2(cd + fg)\sigma_3\tau_0 \}^2 + (2ac\sigma_1\tau_0 + 2bd\sigma_1\tau_3)^2 (2ag\sigma_2\tau_3 + 2bf\sigma_2\tau_0)^2$$

$$= 4\{(ab)^2 + (cd + fg)^2 + (ac)^2 + (bd)^2 + (ag)^2 + (bf)^2 \} - 8ab(cd + fg)\sigma_0\tau_3 + 8acbd\sigma_0\tau_3 + 8agbf\sigma_0\tau_3$$

$$= 4\{(ab)^2 + (cd + fg)^2 + (ac)^2 + (bd)^2 + (ag)^2 + (bf)^2 \}.$$
(B15)

Therefore, the eigenvalues are written as

$$E'_{\rm p\pm} = x \pm \sqrt{y^2 + C'_{\rm p}^2},\tag{B16}$$

$$E'_{\rm m\pm} = x \pm \sqrt{y^2 + C'_{\rm m}^2},\tag{B17}$$

with

$$C_{\rm p}^{\prime 2} = a^2 + b^2 + c^2 + d^2 + f^2 + g^2 + 2\sqrt{(ab)^2 + (cd + fg)^2 + (ac)^2 + (bd)^2 + (ag)^2 + (bf)^2},$$
 (B18)

$$C_{\rm m}^{\prime 2} = a^2 + b^2 + c^2 + d^2 + f^2 + g^2 - 2\sqrt{(ab)^2 + (cd + fg)^2 + (ac)^2 + (bd)^2 + (ag)^2 + (bf)^2}.$$
 (B19)

Thus, choosing the parameters as Eq. (B13), we obtain Eq. (B10).

3. Numerical results

In the main text, we have briefly discussed the extended Hatano-Nelson chain. Here, let us numerically analyze this system in detail. Firstly, we focus on the subsector with (N, P) = (3, -1) [see Fig. 6]. Although the topology of the one-body Hamiltonian is non-trivial [i.e., $(w, w_s) = (0, 1)$ for $\epsilon_{ref} = 0$], the many-body Hamiltonian is topologically trivial [i.e., $W_{(3,-1)} = 0$ for $E_{ref} = 0$] as shown in Figs. 6(a) and 6(b). This fact results in the fragility of the non-Hermitian skin effect against interactions. Namely, although the fermion with the up- (down-) spin state is localized at the right (left) edge due to the non-Hermitian skin effect in the noninteracting case [see Figs. 6(c) and 6(e)], such localization cannot be observed in the presence of the interactions [see Figs. 6(d) and 6(f)]. Correspondingly, the extreme sensitivity of the energy spectrum to the boundary condition is not observed for J = V = 1 [see Figs. 3(b) and 3(d)].

This fragility of the non-Hermitian skin effect is intuitively understood as follows: the interactions flip the spin of fermions in orbital a, which suppresses the effects of the boundaries.

Now, let us focus on the subsector with (N, P) = (4, 1)[see Figs. 7 and 8]. Figures 7(a) and 7(b) indicate the topology of the many-body Hamiltonian is trivial. How-



FIG. 6. Data for (L,t) = (7,1) and the subsector (N,P) = (3,-1). (a) and (b): The twist angle (θ) dependence of $\arg[\prod_n (E_n - E_{ref})]$ for $E_{ref} = 0$. (c) and (d) [(e) and (f)]: Expectation values $\langle \hat{n}_{ja\sigma} \rangle = {}_{\mathrm{R}} \langle \Phi_n | \hat{n}_{ja\sigma} | \Phi_n \rangle_{\mathrm{R}}$ with $\sigma =\uparrow [\sigma =\downarrow]$. Here, $|\Phi_n\rangle_{\mathrm{R}}$ (n = 0, 1, ...) denote right eigenstates of \hat{H}_{eHN} . Data obtained under the open boundary condition (the periodic boundary condition) are shown with colored (gray) symbols. Panels (a), (c), and (e) [(b), (d), and (f)] display data for J = V = 0 [J = V = 1].

ever, due to the topology of the one-body Hamiltonian, we can observe the extreme sensitivity of the spectrum and expectation values $\langle \hat{n}_{ja\sigma} \rangle$ to the presence/absence of the boundaries [see Figs. 7(c) and 7(d)]. As is the case for (N, P) = (3, -1), such extreme sensitivity is fragile against interactions [see Fig. 8].

Finally, we discuss the case for (N, P) = (9, -1) where orbital *a* is half-filled. Figure 9 indicates that the manybody Hamiltonian is topologically trivial, which results in fragility of the non-Hermitian skin effect at the noninteracting level as discussed in the above. Namely, while the topology of the one-body Hamiltonian induces the extreme sensitivity of the energy spectrum and the expectation values $\langle \hat{n}_{ja\downarrow} \rangle$ to the boundary conditions, such extreme sensitivity is not observed in the interacting case [see Fig. 10].



FIG. 7. Numerical data for J = V = 0 and the subsector (N, P) = (4, 1). (a): Spectral flow of the manybody Hamiltonian. (b): The twist angle (θ) dependence of $\arg[\prod_n (E_n - E_{ref})]$ for $E_{ref} = 0.3i$. (c) and (d) [(e) and (f)]: Expectation values $\langle \hat{n}_{ja\sigma} \rangle = {}_{\rm R} \langle \Phi_n | \hat{n}_{ja\sigma} | \Phi_n \rangle_{\rm R}$ with $\sigma = \uparrow [\sigma = \downarrow]$. Here, $|\Phi_n \rangle_{\rm R}$ (n = 0, 1, ...) denote right eigenstates of $\hat{H}_{\rm eHN}(\theta = 0)$. Data obtained under the open boundary condition (the periodic boundary condition) are shown with colored (gray) symbols. These data are obtained for a parameter set (L, t) = (7, 1).



FIG. 8. Numerical data for J = V = 1 and the subsector (N, P) = (4, 1). These figures are plotted in the same way as Figs. 7(a)-7(d).



FIG. 9. Data for (L,t) = (7,1) and the subsector (N,P) = (9,-1). (a) and (b) [(c) and (d)]: Spectral flow for J = V = 0 [J = V = 1]. Panel (b) [(d)] is a magnified version of the range $-0.1 \leq \text{Im}E \leq 0.1$ and $-0.1 \leq \text{Re}E \leq 0.1$ in panel (a) [(c)]. (e) and (f): The twist angle (θ) dependence of $\arg[\prod_n (E_n - E_{\text{ref}})]$ for $E_{\text{ref}} = -0.04$ and J = V = 0 [J = V = 1].



FIG. 10. (a) and (b): Spectrum of the many-body Hamiltonian for J = V = 0 and J = V = 1, respectively. Blue (red) symbols denote data obtained under the periodic (open) boundary condition. (c) and (d): Expectation values $\langle \hat{n}_{ja\downarrow} \rangle = {}_{\mathrm{R}} \langle \Phi_n | \hat{n}_{ja\downarrow} | \Phi_n \rangle_{\mathrm{R}}$ against j and $\langle \hat{N}_{a\uparrow} \rangle = {}_{\mathrm{R}} \langle \Phi_n | \hat{N}_{a\uparrow} | \Phi_n \rangle_{\mathrm{R}}$ for J = V = 0 and J = V = 1, respectively. These data are obtained for (L, t) = (7, 1) and the subsector (N, P) = (9, -1).