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Hardware-Encoding Grid States in a Nonreciprocal Superconducting Circuit

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We present a circuit design composed of two Josephson junctions coupled by a nonreciprocal element, the gyrator, whose ground space is doubly degenerate. The ground states are approximate code words of the Gottesman-Kitaev-Preskill code. We determine the low-energy dynamics of the circuit by working out the equivalence of this system to the problem of a single electron in a crystal, confined to a two-dimensional plane, and subjected to a strong, homogeneous magnetic field. We find that the circuit is naturally protected against the common noise channels in superconducting circuits, such as charge and flux noise, implying that it can be used for passive quantum error correction. We also propose realistic design parameters for an experimental realization, and we describe possible protocols to perform logical one- and two-qubit gates, state preparation, and readout.

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I. INTRODUCTION

Building a quantum computer in a physical system is a formidably challenging task because of the inherent fragility of physical quantum bits (qubits). The key idea behind quantum error correction (QEC) [1,2] is to use *logical* qubits that can be protected against certain likely errors, thus extending the lifetime of the encoded quantum information and allowing for fault-tolerant quantum computation [3,4].

There are different flavors of QEC codes, that differ in the way in which the logical qubits are constructed. For example, in the toric [5,6], surface [7–9], and color [10] code, the logical qubits are encoded in lattices of physical qubits. To date, the QEC codes that have been most successful in enhancing the lifetime of quantum information have been built from continuous variable (CV) systems [11–13], such as a single microwave cavity mode. Efficient QEC with cat states and binomial codes has been demonstrated [14,15]. In this work, we focus on a similar CV encoding, proposed by Gottesman, Kitaev, and Preskill (GKP) in Ref. [16], where the code words are shifted grid states, i.e., periodic repetitions of highly localized peaks

such that the code words do not have common support. With the GKP code, these code words can be protected against sufficiently small displacements in phase space. The error-correcting properties of the GKP code are further explored in Ref. [17], where it is shown that the GKP code outperforms cat and binomial codes when a photon loss channel is considered. Grid states have been successfully prepared and actively stabilized in a superconducting microwave cavity by applying conditional displacements depending on the measurement outcome of an ancillary physical qubit [18]. Also, implementations of these states in superconducting circuits have been proposed by means of the $0-\pi$ qubit [19,20] or in a passive procedure with the dualmon [21].

However, the *active* implementation of QEC requires complicated protocols where errors are detected and compensated for by applying a recovery operation. In contrast, in *passive* QEC, the protection is a built-in feature of the system's hardware, and it is, therefore, advantageous in terms of hardware efficiency and scalability. Generally, this is achieved by constructing a system whose twofold degenerate ground states are the qubit states: Errors that bring the system out of the computational space have an associated energy penalty, and so the system automatically relaxes back into the computational space [5,22,23].

An example of an implementation of the GKP code is a single electron confined in a two-dimensional plane within a periodic potential and a high perpendicular magnetic field [24–37]. The magnetic field restricts the dynamics of the electron to the lowest Landau level (LLL),

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so that the position operators in orthogonal directions do not commute. The contribution of the periodic potential to the Hamiltonian reduces to a sum of displacement operators, which are the stabilizers of the GKP code.

Although this system is useful for a theoretical understanding of the code, it is very unpractical to implement. The magnetic field required for it to work is exactly $B = \Phi_0/2A$ with A the area of the unit cell in space and the magnetic flux quantum $\Phi_0 = h/e$. In natural crystals, this condition implies that the magnetic field needs to be unrealistically large, about $B \sim 10^5$ T. Moiré patterns in twisted bilayer graphene can be used to reduce this value by a few orders of magnitude due to their large unit cell [38]. Even if such a regime were possible to achieve, the external magnetic field would still require an extremely precise fine-tuning, and the electron density would still have to be decreased to the unfeasible value of a single free electron in the crystal.

Here, we propose instead a different implementation of the GKP code, which does not suffer from any of these issues. We consider a superconducting circuit composed of two Josephson junctions coupled by a lossless and linear nonreciprocal circuit element, the *gyrator* [39]. The circuit is shown in Fig. 1.

The nonreciprocity of the gyrator breaks time-reversal symmetry, and its contribution to the dynamics of the circuit is akin to that of a uniform, perpendicular magnetic field in an electronic system. In our circuit, the condition on the strength of the magnetic field reduces to the requirement on the gyration conductance being precisely twice the conductance quantum, i.e., $G = 2e^2/h$.

While unrealistic for conventional gyrators [40–45], this value of G can be easily reached in quantum Hall effect (QHE) devices [46–50], in which a precise fine-tuning of the device parameters is unnecessary due to the quantization of the off-diagonal conductivity. In addition, although conventional QHE devices [50] require a high magnetic field to operate, making them unpractical to couple to superconducting devices, we note that state-of-the-art quantum anomalous Hall effect materials still present an extremely precise conductivity quantization and low losses [51] and can be used to implement nonreciprocal electrical network elements that can operate at zero magnetic field [52].

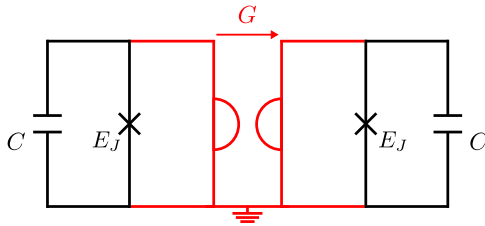


FIG. 1. Proposed hardware implementation of the GKP code. The circuit consists of two Josephson junctions coupled by a gyrator, highlighted in red.

We show that our construction is insensitive to common types of noise. We discuss possible ideas of how logical one-qubit as well as two-qubit Clifford gates can be implemented by applying currents and using tunable inductances. We show that the ground state of our system is an eigenstate of the Hadamard gate. Consequently, our system is suitable for universal quantum computation [16,53,54].

The paper is organized as follows: In Sec. II, we review a few key concepts of hardware-encoding GKP states, and we introduce the Hamiltonian whose twofold degenerate ground space is spanned by the GKP code words. In Sec. III, we show how this Hamiltonian can be derived from the low-energy description of a single electron in a high magnetic field and a periodic potential. The dynamics of this system is equivalent to that of a gyrator connected to two Josephson junctions, but the solid-state jargon more easily reveals the intimate relation to Hofstadter’s butterfly [34,37]: The GKP states are obtained at a specific point in the butterfly. In Sec. IV, we study the effect of an additional parabolic confinement potential, which in the circuit model consists of the addition of inductances in parallel to the Josephson junctions. For this setting, the ground space of the resulting Hamiltonian is twofold degenerate up to an exponentially small gap, and the eigenstates of the system resemble superpositions of normalizable, approximate GKP code words [16]. In Sec. V, we highlight the connection between the one-dimensional GKP grid states and the two-dimensional ground space wave functions of the electronic system’s Hamiltonian projected onto the lowest Landau level. In Sec. VI, we work out in detail the equivalence to the nonreciprocal superconducting circuit model, and we propose realistic design parameters for an experimental realization of the system. We also discuss possible realizations of logical gates by using current sources and tunable inductances, and we present ideas for state preparation and readout. We provide an analysis of the protection against common noise sources, such as flux and charge noise. In Sec. VII, we compare our qubit design with other currently existing or proposed superconducting qubits, and we discuss possible trade-offs of our design. Finally, in Sec. VIII, we summarize our results and give an outlook on further work.

II. THE GKP CODE FOR PASSIVE QEC

The GKP code is a CV quantum error-correcting code [11–13] introduced in Ref. [16]. In contrast to the standard approach to quantum error correction, which assumes physical qubits as fundamental noisy elements, in CV quantum error correction, the idea is to encode a two-level (or n -level, more generally) system in the infinite-dimensional Hilbert space of a one-dimensional particle characterized by dimensionless canonical quadrature operators X and P satisfying $[X, P] = i$. The GKP code can be described within the stabilizer formalism for CV systems.

The role of the Pauli group is played by the Weyl-Heisenberg group G_H , i.e., the group of displacement operators [55,56]

$$D(\alpha) = e^{ab^\dagger - \alpha^* b}, \quad \alpha \in \mathbb{C}, \quad (1)$$

with the annihilation operator $b = (X + iP)/\sqrt{2}$. In this framework, the GKP code is the two-dimensional subspace stabilized by a subgroup $\mathcal{S}_{\text{GKP}}$ of G_H with group generators

$$S_X = D(i\sqrt{2\pi}) = e^{i2\sqrt{\pi}X}, \quad S_P = D(\sqrt{2\pi}) = e^{-i2\sqrt{\pi}P} \quad (2)$$

and S_X^{-1}, S_P^{-1} . The logical Pauli operators $\bar{Z}$ and $\bar{X}$ are given by

$$\bar{Z} = S_X^{1/2} = e^{i\sqrt{\pi}X}, \quad \bar{X} = S_P^{1/2} = e^{-i\sqrt{\pi}P}. \quad (3)$$

This choice of logical operators fixes the following (unnormalizable) code words:

$$|\bar{0}\rangle = \sum_{n \in \mathbb{Z}} |X = 2\sqrt{\pi}n\rangle, \quad (4a)$$

$$|\bar{1}\rangle = \sum_{n \in \mathbb{Z}} |X = 2\sqrt{\pi}n + \sqrt{\pi}\rangle, \quad (4b)$$

that are grid states, each describing a comb of equidistant δ peaks in the X basis. These combs have a period of $2\sqrt{\pi}$ and are shifted with respect to each other by $\sqrt{\pi}$.

Given a density matrix ρ describing the state of a one-dimensional CV quantum system, one can expand a generic quantum operation $\mathcal{E}(\rho)$ in terms of displacement operators. For simplicity, we consider the superoperator of a diagonal channel, which can be expressed as [16]

$$\mathcal{E}(\rho) = \int_{\mathbb{C}} d\alpha f(\alpha) D(\alpha) \rho D^\dagger(\alpha), \quad (5)$$

with $f(\alpha)$ being a scalar function. If $D^\dagger(\alpha')D(\alpha)$ either is in the stabilizer group or does not commute with the stabilizers S_X and S_P in Eq. (2), the GKP code featuring the ideal code words in Eq. (4) can correct against these kinds of errors. Mathematically, this condition means that $f(\alpha)$ must have support only for

$$|\text{Re}(\alpha)| < \frac{1}{2}\sqrt{\frac{\pi}{2}}, \quad |\text{Im}(\alpha)| < \frac{1}{2}\sqrt{\frac{\pi}{2}}. \quad (6)$$

In this case, the error syndromes are unique, and the displacement errors $D(\alpha)$ can be corrected. Otherwise, logical errors will be made.

The main idea behind passive, stabilizer error correction is to construct a Hamiltonian that has the code subspace as the low-energy subspace. For the GKP code, this Hamiltonian is easily obtained as [16]

$$H_{\text{GKP}}/V_0 = -[\cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)], \quad (7)$$

with V_0 a constant with the unit of energy. Because the code subspace is stabilized by the two cosines, it has energy $-2V_0$.

We remark that passive, stabilizer error correction for CV systems is rather different than in systems based on a large set of physical qubits, such as the toric, surface, or color code [5–10]. In fact, the Hamiltonian in Eq. (7) is gapless, with a continuous spectrum ranging from $-2V_0$ to $+2V_0$. Also, because the Weyl-Heisenberg group is a continuous group, in contrast to the discrete Pauli group, the eigenstates of H_{GKP} are unnormalizable and formally out of the Hilbert space of any physical system. As a consequence, the usual perturbation theory argument [5,23] which claims that *local* perturbations of the Hamiltonian give rise to small variations of the energy levels is not directly applicable here. As pointed out in Ref. [22], the argument can be restored in approximated versions of Eq. (7), where the eigenstates are normalized and confined, leading to a discrete spectrum. The particular way in which H_{GKP} is approximated modifies the properties of the degenerate ground space, but generally its eigenstates remain with disjoint support. This idea of passive protection is similar to the one behind the 0- π qubit [19,20,57–61].

Finally, the addition of a confinement to H_{GKP} provides ways to prepare approximate versions of the GKP code words in Eq. (4), or linear combinations of them, which become the quasidegenerate ground states of the confined system. However, for the GKP code featuring approximate grid states, the condition of an error being correctable is more intricate than Eq. (6) [16].

III. CRYSTAL ELECTRON IN A MAGNETIC FIELD

We discuss how the code Hamiltonian in Eq. (7) can emerge from the consideration of the well-known situation of a single electron confined to a two-dimensional plane in a strong perpendicular uniform magnetic field [24–37]. The effect of a periodic potential on the electron's motion has been extensively studied because of the fractal nature of the energy bands [34] and their nontrivial topology [62]. In this section, we want to clarify under what conditions the ground states of the system are GKP states. We focus on the Hamiltonian

$$H = \frac{[\mathbf{p} + e\mathbf{A}(x_1, x_2)]^2}{2m} + V_{\text{crys}}(x_1, x_2), \quad (8)$$

where the two-dimensional positions and momenta satisfy canonical commutation relations $[x_i, p_j] = i\hbar\delta_{ij}$, for $i, j \in \{1, 2\}$, and $\mathbf{B} = \nabla \times \mathbf{A} = B\mathbf{e}_3$. We consider a crystal potential of the form

$$V_{\text{crys}}(x_1, x_2) = -V \left[\cos\left(2\pi \frac{x_1}{L_0}\right) + \cos\left(2\pi \frac{x_2}{L_0}\right) \right], \quad (9)$$

which corresponds to the first Fourier mode of any periodic potential on a square lattice of size L_0 in the $x_1 x_2$ plane. Although both the crystal potential V_{crys} and the uniform magnetic field $\mathbf{B}$ are periodic in the x_1 and x_2 direction, the Hamiltonian is not, because the discrete translation symmetry is broken in at least one direction by the vector potential $\mathbf{A}$. As a result, H does not simultaneously commute with both the canonical unitary translation operators $t_1(L_0)$ and $t_2(L_0)$, defined as

$$t_i(r) = e^{-irp_i/\hbar}, \quad r \in \mathbb{R}, \quad i = 1, 2. \quad (10)$$

It follows that H , $t_1(L_0)$ and $t_2(L_0)$ cannot have common eigenstates, as the usual formulation of Bloch's theorem dictates. To find a set of translations which do commute with the Hamiltonian, we work with the dynamical momenta [63–66]

$$\pi_1 = p_1 + eA_1, \quad \pi_2 = p_2 + eA_2 \quad (11)$$

and the guiding center variables

$$R_1 = x_1 - \frac{1}{m\omega_c} \pi_2, \quad R_2 = x_2 + \frac{1}{m\omega_c} \pi_1, \quad (12)$$

with the cyclotron frequency $\omega_c = eB/m$. These operators are gauge invariant (in contrast to the canonical momenta p_i) and satisfy the commutation relations

$$[\pi_1, \pi_2] = -i \frac{\hbar^2}{l_B^2}, \quad [R_1, R_2] = i l_B^2, \quad [\pi_i, R_j] = 0, \quad (13)$$

for $i, j \in \{1, 2\}$ and the magnetic length $l_B = \sqrt{\hbar/eB}$. Physically, the dynamical momenta are related to the cyclotron motion of an electron around its center of mass, which, in turn, is parametrized by the guiding center coordinates.

We can impose boundary conditions by requiring the wave function to be quasiperiodic in R_i . To this end, we make use of the unitary *magnetic* translation operators (MTOs) [27,66–69]

$$T_1(r) = e^{-irR_2/l_B^2}, \quad T_2(r) = e^{irR_1/l_B^2}, \quad r \in \mathbb{R}, \quad (14)$$

which shift the guiding center variables R_i by r , i.e.,

$$T_i^\dagger(r) R_i T_i(r) = R_i + r, \quad i = 1, 2. \quad (15)$$

It is straightforward to show that the MTOs $T_1(L_0)$ and $T_2(L_0)$ do commute with the Hamiltonian in Eq. (8).

However, because of the noncommutativity of R_1 and R_2 in Eq. (12), magnetic translations in different directions do not generally commute. An electron moving along a closed path accumulates an Aharonov-Bohm phase [70] proportional to the magnetic flux threaded by the loop, and so

$$T_2(r_2) T_1(r_1) = e^{i2\pi B r_1 r_2 / \Phi_0} T_1(r_1) T_2(r_2), \quad (16)$$

with the (nonsuperconducting) flux quantum $\Phi_0 = h/e$. Consequently, MTOs in orthogonal directions commute only when an integer number of flux quanta is threaded through the loop defined by the MTOs. Note that, with an appropriate rescaling, the MTOs defined in Eq. (14) correspond to the displacement operators similar to the ones defined in Eq. (1).

In the following, we restrict to *rational* fluxes [34], where the magnetic flux enclosed in a unit cell of size $L_0 \times L_0$ is a rational multiple of the flux quantum, i.e.,

$$\Phi = BL_0^2 = \frac{p}{q} \Phi_0, \quad (17)$$

with coprime natural numbers p and q . In this case, we consider an enlarged, *magnetic* unit cell of size $qL_0 \times L_0$, which contains p flux quanta, such that the MTOs $T_1(qL_0) = [T_1(L_0)]^q$ and $T_2(L_0)$ commute with the Hamiltonian in Eq. (8) and with each other [71]. As a result, we consider the magnetic Bloch states satisfying

$$T_1(qL_0)|\mathbf{k}\rangle = e^{ik_1 qL_0} |\mathbf{k}\rangle, \quad (18a)$$

$$T_2(L_0)|\mathbf{k}\rangle = e^{ik_2 L_0} |\mathbf{k}\rangle, \quad (18b)$$

where $\mathbf{k} = (k_1, k_2)^T$ is the crystal momentum defined in the rectangular Brillouin zone

$$k_1 \in \left[0, \frac{2\pi}{qL_0}\right), \quad k_2 \in \left[0, \frac{2\pi}{L_0}\right). \quad (19)$$

The states $|\mathbf{k}\rangle$ are sometimes referred to as Zak states [28–30].

Introducing the Landau-level ladder operators

$$a = \frac{1}{\sqrt{2}} \frac{l_B}{\hbar} (\pi_2 + i\pi_1), \quad a^\dagger = \frac{1}{\sqrt{2}} \frac{l_B}{\hbar} (\pi_2 - i\pi_1), \quad (20)$$

satisfying the bosonic commutation relation $[a, a^\dagger] = 1$, the Hamiltonian in Eq. (8) can be rewritten as

$$H = \hbar\omega_c \left(a^\dagger a + \frac{1}{2} \right) - \frac{V}{2} \left[D_a \left(\frac{i\sqrt{2}\pi l_B}{L_0} \right) T_1 \left(\frac{qL_0}{p} \right) + D_a \left(\frac{-\sqrt{2}\pi l_B}{L_0} \right) T_2 \left(\frac{qL_0}{p} \right) + \text{H.c.} \right], \quad (21)$$

with the cyclotron frequency $\omega_c = eB/m$ and the unitary displacement operator $D_a(\alpha) = \exp(\alpha a^\dagger - \alpha^* a)$ acting on the subspace of the dynamical momenta. A convenient basis to numerically analyze the low-energy spectrum of this Hamiltonian are the product states $|n; \mathbf{k}, l\rangle = |n\rangle \otimes |\mathbf{k}, l\rangle$ satisfying

$$a^\dagger a |n\rangle = n |n\rangle, \quad n \in \mathbb{N}_0, \quad (22)$$

and

$$T_1(qL_0) |\mathbf{k}, l\rangle = e^{ik_1 q L_0} |\mathbf{k}, l\rangle, \quad (23a)$$

$$T_2(L_0/p) |\mathbf{k}, l\rangle = e^{i(k_2 L_0 + 2\pi l)/p} |\mathbf{k}, l\rangle, \quad (23b)$$

$$\langle \mathbf{k}, l | \mathbf{k}', l' \rangle \propto \delta_{ll'}, \quad (23c)$$

with $l = 0, 1, 2, \dots, p-1$; see the Appendix A. Note that, in the absence of the potential ($V = 0$), the states $|n; \mathbf{k}, l\rangle$ diagonalize the Hamiltonian, leading to a p -fold degeneracy of each Landau level with energy $E_n = \hbar\omega_c(n + 1/2)$ that does not depend on $\mathbf{k}$. This degeneracy is the well-known flat Landau-level degeneracy, which is often discussed in the quantum Hall effect literature [63–65]. Expressed in the $x_1 x_2$ representation, the quasiperiodic wave functions $\Psi_{n;\mathbf{k},l}(x_1, x_2) = \langle x_1, x_2 | n; \mathbf{k}, l \rangle$ are introduced by Haldane and Rezayi in Ref. [72]; see also Sec. V.

The crystal potential couples states with different Landau-level occupation numbers n and with different guiding center quantum numbers l . Consequently, the p -fold degeneracy of each initially flat Landau level is lifted, and, for a moderately weak crystal potential, each Landau level splits into p subbands with finite broadening [31–33]; see Fig. 2, in which the two lowest split Landau

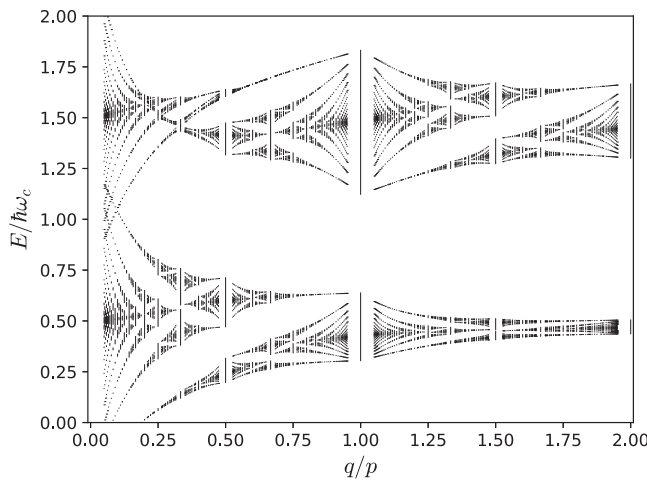


FIG. 2. Low-energy spectrum of the Hamiltonian in Eq. (21) as a function of the *inverse* flux ratio q/p for a fixed value of $V/\hbar\omega_c = 0.4$. The initially flat Landau levels at $E_n/\hbar\omega_c = n + 1/2$ (obtained for $V = 0$) are split into subbands dependent on $\mathbf{k}$.

levels are shown. Moreover, although each subband has a functional dependency on $\mathbf{k}$, each band has an additional q -fold degeneracy; see Appendix A. More details on the general solution of the Hamiltonian in Eq. (21) can be found in Appendix B 1.

In this paper, we are interested in the weak Landau-level coupling limit $V/\hbar\omega_c \ll 1$, where the dynamics of states within each Landau level can be taken to be independent from the others. This limit is analyzed in the following.

A. GKP qubit in the LLL projection

When the coupling between the Landau levels is weak, an effective low-energy Hamiltonian acting on a single Landau level can be obtained by a Schrieffer-Wolff transformation [73,74]. In particular, to the lowest order and considering only the LLL [75], we obtain from Eq. (21) the effective Hamiltonian (up to an unimportant constant)

$$H_{\text{LLL}} = \langle n=0 | H | n=0 \rangle = -\frac{V_0}{2} \left[T_1\left(\frac{qL_0}{p}\right) + T_2\left(\frac{qL_0}{p}\right) + \text{H.c.} \right], \quad (24)$$

where $V_0 = V e^{-\pi q/2p}$. Although formally this effective Hamiltonian is valid only when $V/\hbar\omega_c \ll 1$, numerics shows that the approximation holds up well to relatively high values of $V/\hbar\omega_c \lesssim 0.4$; see Sec. IV A, where we discuss in more detail the validity of the LLL projection.

In the limit of weak Landau-level coupling, the eigenvalue equation associated with H_{LLL} is the Harper equation [25,26,35,37,76], which is a special case of the almost Mathieu equation [77–79]. In particular, the Harper equation is a finite-difference equation, resulting in an energy spectrum in form of the Hofstadter butterfly [34], shown in Fig. 3.

We point out that this spectrum has p bands that are q -fold degenerate. In fact, states connected by the

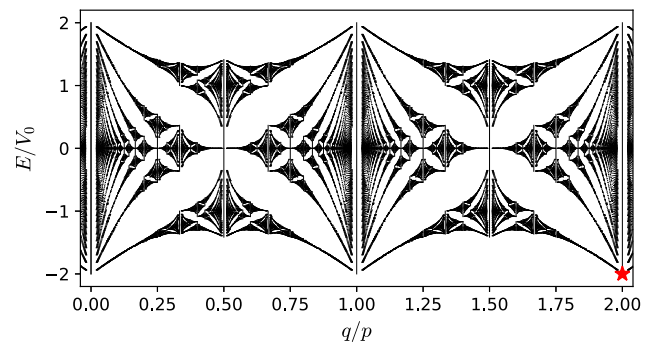


FIG. 3. Hofstadter's butterfly obtained by plotting the spectrum of the effective lowest Landau-level Hamiltonian H_{LLL}/V_0 , defined in Eq. (24), as function of the *inverse* magnetic flux ratio $\Phi_0/\Phi = q/p$. The two-dimensional GKP code space corresponds to the states of minimal energy at $p/q = 1/2$. This point is marked by a red star in the figure.

application of $T_1(nL_0)$ with $n = 1, \dots, q-1$ are orthogonal but have the same energy; see Appendix A. Note that this result is in contrast to the original tight-binding result of Hofstadter [34], where different Landau levels are strongly coupled and where there are q bands that are q -fold degenerate [76]. For this reason, in the original work, the Hofstadter butterfly is obtained by plotting the spectrum as a function of p/q instead of q/p [34].

Importantly, by introducing the dimensionless variables

$$X = \frac{\sqrt{\pi}}{L_0} R_1 = \sqrt{\frac{q}{2p}} \frac{R_1}{l_B}, \quad P = \frac{2p\sqrt{\pi}}{qL_0} R_2 = \sqrt{\frac{2p}{q}} \frac{R_2}{l_B}, \quad (25)$$

satisfying the canonical commutation relation $[X, P] = i$, we can rewrite H_{LLL} as

$$H_{\text{LLL}}/V_0 = -\left[\cos(2\sqrt{\pi}X) + \cos\left(\frac{q}{p}\sqrt{\pi}P\right)\right]. \quad (26)$$

Comparing with Eq. (7), we observe that H_{LLL} corresponds to the GKP Hamiltonian H_{GKP} when $p/q = 1/2$. This system is, therefore, suitable for passively encoding the GKP code words, which are given in Eq. (4). Note that, for $p/q = 1/2$, the previous rescaling of the guiding center variables R_i becomes equal, and so the code can correct equal shifts on X and P .

Furthermore, for $p/q = 1/2$, the MTOs defining the states $|k\rangle$ in Eq. (18) are related to the stabilizers and logical operators of the GKP code as, respectively,

$$T_1(2L_0) = S_P, \quad T_2(L_0) = S_X^{1/2} = \bar{Z}. \quad (27)$$

Since the GKP code words are the eigenstates of H_{LLL} with minimal eigenenergy, we can identify the code space with a specific point in Hofstadter's butterfly (see the red star in Fig. 3). In particular, the code space is spanned by the eigenstates obtained for $k = (0, 0)^T$ and $k = (0, \pi/L_0)^T$. These states correspond to the logical code words introduced in Eq. (4). The eigenfunctions of the full system within the LLL projection are analyzed in Sec. V.

At this point, we highlight the main difference between our approach and the original proposal by GKP [16]. GKP propose to use the LLL projection at the rational flux $p/q = d/1$ without the crystal potential ($V_{\text{crys}} \equiv 0$). A qudit can be encoded by focusing on the d -fold degenerate ground space obtained at vanishing Bloch momentum ($k = 0$), and one can take this qudit to construct different shift-resistant quantum codes. In contrast, for the rational flux $p/q = 1/2$, by including the crystal potential and using states with different Bloch momenta, here we encode a qubit in a real CV system.

IV. ADDITIONAL PARABOLIC CONFINEMENT

Because the GKP code words in Eq. (4) are not normalizable, they are mathematical objects that are not physically realizable. As a consequence, we can work only with normalized wave packets that approximate the ideal code words. This consideration leads to the introduction of several versions of approximate GKP states [80]. In what follows, we show that the addition of a parabolic confinement allows us to obtain low-energy eigenstates with a finite energy gap that are approximate GKP code words. Thus, the ability to prepare the unique ground state provides a reliable way to initialize the system. Importantly, we show that by including a parabolic confinement potential it is possible to prepare directly approximate GKP magic states, namely, Hadamard eigenstates, which, together with Clifford operations and Pauli measurements, enable universal quantum computation [53].

We start by choosing an isotropic parabolic confining potential, and so the Hamiltonian is

$$H = \frac{[p + eA(x_1, x_2)]^2}{2m} + V_{\text{tot}}(x_1, x_2), \quad (28)$$

with

$$V_{\text{tot}}(x_1, x_2) = V_{\text{crys}}(x_1, x_2) + \frac{1}{2}m\omega_0^2(x_1^2 + x_2^2). \quad (29)$$

Because V_{tot} does not preserve the discrete translational symmetry defined by V_{crys} , we cannot impose the magnetic Bloch conditions in Eq. (18). In this case, we require instead that the wave functions vanish at infinity.

We now briefly summarize the main findings of a numerical analysis of the eigensystem of the Hamiltonian in Eq. (28) (see Appendix B 2 for details). In particular, we focus on the case $p/q = 1/2$ [see Eq. (17)], which in the LLL projection leads to ideal GKP states in the absence of the parabolic confinement potential. As in the previous section, we construct and analyze a low-energy theory of the system, valid in the weak Landau-level coupling limit.

A. Numerical results

When a confinement potential is included, the twofold degeneracy of the ground space is lifted, and an energy gap opens between the degenerate ground states of H_{LLL} in Eq. (26). This energy gap, however, is exponentially small in ω_0 , and so the ground state and the first excited state remain quasidegenerate when ω_0 is small enough. To illustrate this point, in Fig. 4, we show the lowest ten eigenenergies of the Hamiltonian in Eq. (28) as functions of the confinement strength $\hbar\omega_0/V$ for a rather large value of $V/\hbar\omega_c = 0.4$. Recall that V denotes the amplitude of V_{crys} in Eq. (9). We observe that the energy gap between the ground state and first excited state is negligibly small up to confinements of $\hbar\omega_0/V \lesssim 0.2$ and that it nicely fits an

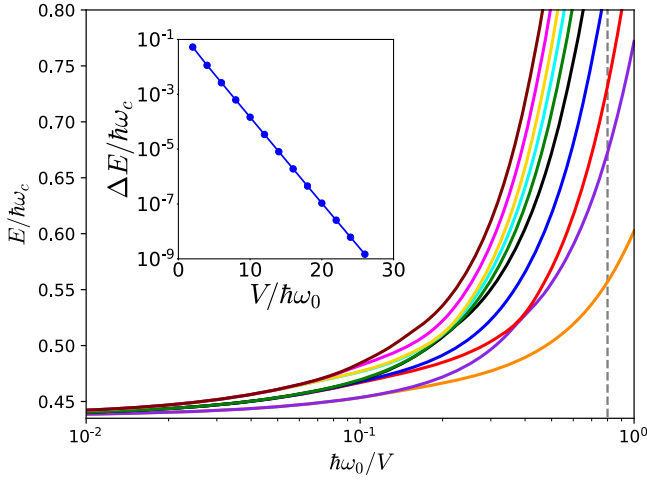


FIG. 4. Lowest ten eigenenergies of the Hamiltonian in Eq. (28) as functions of the confinement strength $\hbar\omega_0/V$ for $p/q = 1/2$ and $V/\hbar\omega_c = 0.4$. The gray dashed vertical line marks the value of $\hbar\omega_0/V = 0.8$ that is experimentally relevant for our circuit proposal; see Sec. VI. Inset: Energy gap $\Delta E = E_1 - E_0$ between the two lowest eigenstates for the same values of p/q and $V/\hbar\omega_c$. For $V/\hbar\omega_0 \gg 1$, the energy gap decreases exponentially with $V/\hbar\omega_0$.

exponential scaling $E_1 - E_0 \propto \exp(-\alpha V/\hbar\omega_0)$, with a positive constant prefactor α [81]. In addition, the spectrum is now discrete, and higher excited states are gapped from the twofold quasidegenerate ground space [81].

As mentioned in Sec. II, the discreteness of the spectrum then allows one to use a perturbative argument which states that local perturbations do not considerably alter the spectrum of the Hamiltonian.

Because we are interested in the weak Landau-level coupling limit, we analyze the effect of higher Landau levels on the low-energy eigenstates numerically. In Fig. 5, we show how the expectation value of the LLL projector $\Pi_{\text{LLL}} = |0\rangle\langle 0|_\pi$ in the ground state of the Hamiltonian in Eq. (28) varies as a function of the confinement potential for fixed values of $V/\hbar\omega_c$. We observe that the higher Landau levels have a negligible effect for a wide range of parameters, giving an error below 3% for rather large values of both $V/\hbar\omega_c$ and $\hbar\omega_0/V$ and, consequently, justifying even in this case a projection onto the LLL. In particular, at the values $V/\hbar\omega_c = 0.4$ and $\hbar\omega_0/V = 0.8$, which are the relevant parameters for the circuit model presented in Sec. VI, we find $\langle\psi_0|\Pi_{\text{LLL}}|\psi_0\rangle = 0.981$; see the green circle in Fig. 5. The expectation values of Π_{LLL} evaluated for the first excited states of the Hamiltonian in Eq. (28) show a similar behavior.

The effective Hamiltonian of the system in this limit is analyzed in the next section.

B. Approximate grid states in the LLL projection

In analogy to Sec. III A, here we find an effective Hamiltonian that captures the behavior of the system in

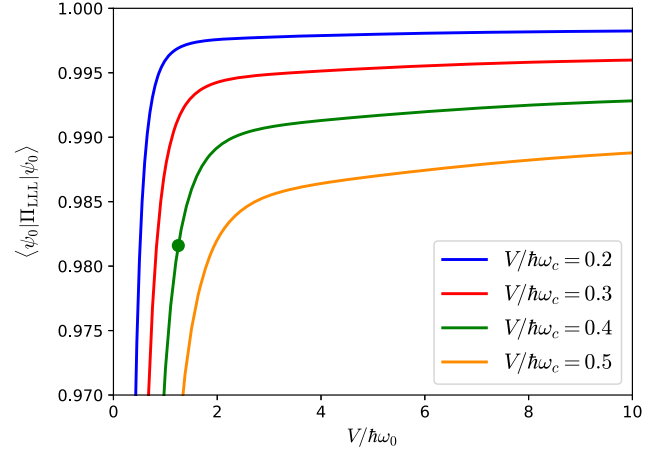


FIG. 5. Expectation value of the LLL projector in the ground state of the Hamiltonian in Eq. (28) for different values of $V/\hbar\omega_c$ as function of $V/\hbar\omega_0$. As expected, the LLL projection becomes more accurate for smaller values of the energy ratios $V/\hbar\omega_c$ and $\hbar\omega_0/V$. We mark with a green circle the values of $V/\hbar\omega_c = 0.4$ and $\hbar\omega_0/V = 0.8$ that are experimentally relevant for our circuit proposal; see Sec. VI. For these values, we find $\langle\psi_0|\Pi_{\text{LLL}}|\psi_0\rangle = 0.981$.

the weak Landau-level coupling limit. By projecting Eq. (28) onto the LLL and considering $p/q = 1/2$, we obtain

$$H_{\text{LLL}} = \frac{\hbar\omega_0^2 P^2 + X^2}{2} - V_0 [\cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)], \quad (30)$$

where $V_0 = Ve^{-\pi}$ and the canonical position X and momentum P are defined by Eq. (25).

Importantly, the parabolic potential in Eq. (29) reduces to the harmonic oscillator Hamiltonian with frequency ω_0^2/ω_c after the LLL projection and breaks the periodicity of both X and P . However, note that the confinement potential preserves the fourfold rotational symmetry in x_1 and x_2 . Because a $\pi/2$ rotation in the x_1x_2 plane reduces to a Fourier transform, which maps $X \mapsto P$ and $P \mapsto -X$, when projected onto the LLL, the Hamiltonian in Eq. (30) is invariant under the exchange of X and P , and its eigenstates are also eigenstates of the Fourier transform. More detailed explanations of this symmetry and correspondence are given in Appendixes B 2 and E.

In particular, the two quasidegenerate low-energy eigenfunctions $\psi_{H+}(X)$ (ground state) and $\psi_{H-}(X)$ (first excited state), shown in Fig. 6, are even and odd eigenfunctions of the Fourier transform with eigenvalues ± 1 , respectively. These states are well approximated by the linear combinations

$$\psi_{H+}(X) \approx \cos\left(\frac{\pi}{8}\right)\psi_0(X) + \sin\left(\frac{\pi}{8}\right)\psi_1(X), \quad (31a)$$

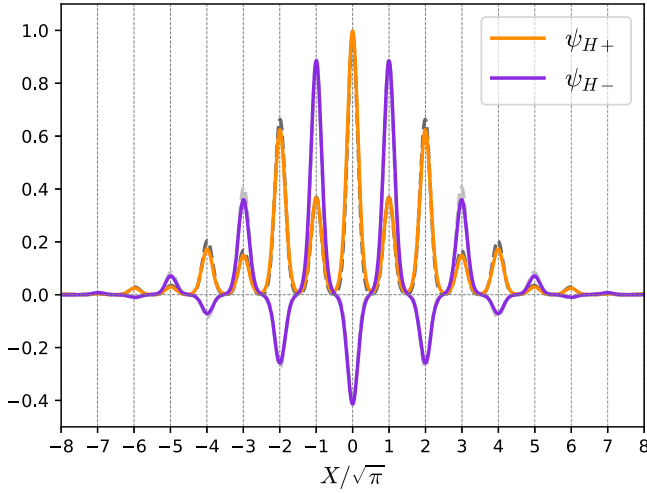


FIG. 6. Numerically obtained lowest-energy eigenfunctions $\psi_{H\pm}(X)$ of the effective Hamiltonian in Eq. (30) for $\hbar\omega_0^2/\omega_c V_0 = 0.05$, which corresponds to $\Delta = 0.25$; see Eq. (33). The quasi-degenerate eigenfunctions are even and odd under Fourier transforming, respectively, and are well approximated by the analytic expressions (dashed light and dark gray lines, respectively) in Eq. (31).

$$\psi_{H-}(X) \approx -\sin\left(\frac{\pi}{8}\right)\psi_0(X) + \cos\left(\frac{\pi}{8}\right)\psi_1(X) \quad (31b)$$

of the approximate grid states

$$\psi_0(X) = \frac{\sqrt{2}}{\pi^{1/4}} e^{-X^2\Delta^2/2} \sum_{n=-\infty}^{\infty} \exp\left(-\frac{(X-2\sqrt{\pi}n)^2}{2\Delta^2}\right), \quad (32a)$$

$$\psi_1(X) = \frac{\sqrt{2}}{\pi^{1/4}} e^{-X^2\Delta^2/2} \sum_{n=-\infty}^{\infty} \exp\left(-\frac{(X-2\sqrt{\pi}n-\sqrt{\pi})^2}{2\Delta^2}\right) \quad (32b)$$

that are shown in Fig. 7. Explicitly, the squeezing parameter Δ is given by

$$\Delta = \left(\frac{\hbar\omega_0^2}{4\pi\omega_c V_0}\right)^{1/4} \ll 1. \quad (33)$$

These approximate grid states are obtained by a convolution of the ideal grid states in Eq. (4) with a narrow Gaussian of width Δ and a multiplication with a wide Gaussian of width $1/\Delta$. This inverse relation is a consequence of the invariance of the Hamiltonian with respect to the Fourier transform, and it also reflects the fact that the GKP code corrects equal errors in the X and P variables.

When the confinement frequency ω_0 is decreased, the width Δ of the individual Gaussian peaks of the approximate grid states in Eq. (32) decreases, while the broadening

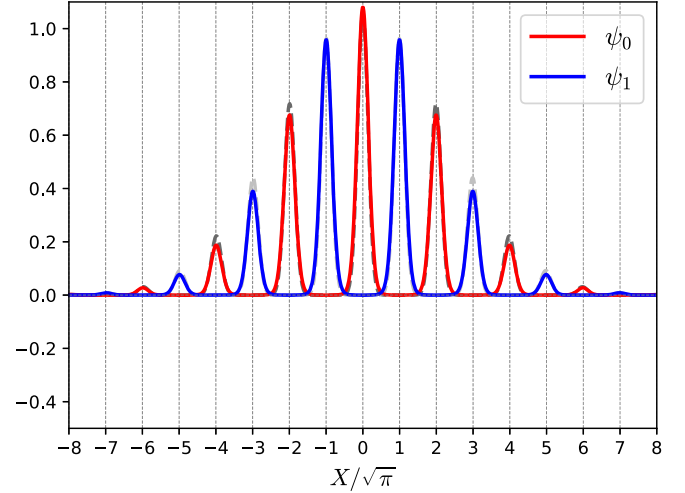


FIG. 7. Approximate GKP code words. We show with solid lines the linear combinations of the numerically obtained lowest-energy eigenfunctions $\psi_{H\pm}(X)$ of Eq. (30) that resemble the approximate GKP states $\psi_{0,1}(X)$; see Eq. (31). The dashed light and dark gray lines show the analytical results in Eq. (32), respectively. The figure is obtained by using the same parameters as in Fig. 6, i.e., $\hbar\omega_0^2/\omega_c V_0 = 0.05$, which corresponds to $\Delta = 0.25$. The widths of the individual peaks and that of the total envelope are the inverse of each other.

of the envelope function increases, eventually recovering the ideal grid states in Eq. (4) when $\omega_0 \rightarrow 0$.

The derivation of Eqs. (31)–(33) is based on a nested application of the envelope function approximation [66,82,83] discussed in Appendix C. Note that, when the parameter Δ is rather small, as we are considering here, the states $\psi_0(X)$ and $\psi_1(X)$ are orthonormal up to an exponentially small correction scaling as approximately e^{-1/Δ^2} . Consequently, they form an appropriate computational basis for the quasidegenerate ground space of H_{LLL} .

The angle $\pi/8$ which appears in the linear combination in Eq. (31) can be understood by considering that, in the basis $\psi_{0,1}(X)$, the Fourier transform approximately equals the Hadamard gate [16]

$$\bar{H} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}, \quad (34)$$

whose even and odd eigenfunctions are $\psi_{H+}(X)$ and $\psi_{H-}(X)$, respectively [84]. These states are magic states which combined with Clifford operations achieve universal quantum computation [16,53,54,85].

The previous result is based on the symmetry of both the amplitudes of the quadratic terms as well as the amplitudes of the cosines in the square brackets in Eq. (30). By breaking this symmetry, for instance, by choosing sufficiently asymmetric amplitudes of the two cosines in Eq. (30), we find that the lowest-energy eigenstates are not Hadamard eigenstates

but the approximate GKP logical code words similar to those shown in Fig. 7. This feature allows one to prepare either Hadamard eigenstates or logical code words by tuning the asymmetry of the crystal potential. We discuss some ideas for state preparation in Sec. VIC.

We remark that an alternative low-energy description of the Hamiltonian in Eq. (28) which relies on the introduction of the eigenbasis of the quadratic part of the Hamiltonian is possible. This basis is known as the Fock-Darwin basis [86,87]. In the weak Landau-level coupling limit, the results obtained with this approach are equivalent to the ones shown here.

V. EIGENFUNCTIONS OF THE TWO-DIMENSIONAL PROBLEM

So far, we describe eigenfunctions of a one-dimensional Hamiltonian obtained by projecting a two-dimensional Hamiltonian onto the LLL. Here, we establish the connection between the eigenfunctions of these two Hamiltonians. In the weak Landau-level coupling regime, which we consider in the previous sections, the wave function of the two-dimensional system is the coherent state representation of the wave function of the one-dimensional system. In the following, we begin by considering the ideal case discussed in Sec. III, and then we straightforwardly generalize our result to include the parabolic confinement potential as introduced in Sec. IV.

In the weak Landau-level coupling limit, the low-energy eigenstates of the Hamiltonian in Eq. (21) are well approximated by the product state

$$|\Psi\rangle = |0\rangle_\pi \otimes |\psi\rangle_R \equiv |0, \psi\rangle, \quad (35)$$

where $|0\rangle_\pi$ denotes the LLL and $|\psi\rangle_R$ is an eigenstate of the projected Hamiltonian in Eq. (26). From Eq. (35), it follows that the wave function $\Psi(x_1, x_2) = \langle x_1, x_2 | \Psi \rangle$ in the original coordinates x_i and the one-dimensional wave function $\psi(X) = \langle X | \psi \rangle_R$ are related by the unitary integral transform

$$\Psi(x_1, x_2) = \int_{-\infty}^{\infty} dX K_0(x_1, x_2; X) \psi(X), \quad (36)$$

where $K_0(x_1, x_2; X) = \langle x_1, x_2 | 0, X \rangle$ is a gauge-dependent integration kernel. As derived in Appendix D, in the symmetric gauge, i.e., $A(x_1, x_2) = B/2(-x_2, x_1, 0)^T$, we obtain

$$K_0(x_1, x_2; X) = \frac{1}{\sqrt{2}\pi^{3/4}} \exp\left(-\frac{(X-x_1)^2}{2}\right) \times \exp(-ix_2 X) \exp\left(i\frac{x_1 x_2}{2}\right). \quad (37)$$

To simplify the notation, in this section, we work in magnetic units, and we rescale all the lengths by the magnetic length $l_B = \sqrt{\hbar/eB}$.

Note that, up to a gauge phase $\exp(ix_1 x_2/2)$, the integration kernel $K_0(x_1, x_2; X)$ is the complex conjugate of the wave function (in X representation) of a coherent state with average position and momentum x_1 and x_2 , respectively. Consequently, as long as the matrix elements between different Landau levels are small and the approximate factorization in Eq. (35) is valid, the low-energy two-dimensional eigenfunctions of the Hamiltonian in Eq. (21) are the coherent state representations of the eigenfunctions of the projected Hamiltonian in Eq. (26) [88]. It follows that the absolute value squared $|\Psi(x_1, x_2)|^2$ is the non-negative Husimi Q representation [56,89,90] associated with $\psi(X)$. Note also that the integral transform in Eq. (36) is invertible and preserves orthonormality.

Let us consider now the LLL non-normalizable eigenfunction

$$\psi_k(X) = e^{-ik_1 X} \sum_{n \in \mathbb{Z}} \delta(X - 2\sqrt{\pi}n - k_2) \quad (38)$$

of the effective Hamiltonian in Eq. (26), satisfying the quasiperiodic boundary conditions defined by Eq. (23) with $p/q = 1/2$ and $l = 0$. Because of the choice $p/q = 1/2$, the length L_0 (in magnetic units) reduces to $\sqrt{\pi}$; see Eq. (17). The wave function of the two-dimensional system, obtained via the integral transform in Eq. (36), is

$$\Psi_k(x_1, x_2) = \frac{1}{\sqrt{2}\pi^{3/4}} e^{-x_1(x_1 - ix_2)/2} \times \vartheta \left[\begin{matrix} k_2/2\sqrt{\pi} \\ -k_1/\sqrt{\pi} \end{matrix} \right] \left(\frac{-i(x_1 - ix_2)}{\sqrt{\pi}}, 2i \right), \quad (39)$$

with the generalized elliptic theta function

$$\vartheta \left[\begin{matrix} a \\ b \end{matrix} \right] (z, \tau) = \sum_{n \in \mathbb{Z}} e^{i\pi(n+a)^2\tau} e^{i2\pi(n+a)(z+b)}. \quad (40)$$

Note that the absolute value of the wave function in Eq. (39) is periodic in both x_1 and x_2 , i.e., $|\Psi_k(x_1, x_2)| = |\Psi_k(x_1 + 2\sqrt{\pi}, x_2)| = |\Psi_k(x_1, x_2 + \sqrt{\pi})|$. Similar two-dimensional functions satisfying quasiperiodic boundary conditions are introduced by Haldane and Rezayi [72] as a basis to describe the problem of an electron confined to the surface of a torus and under the effect of a perpendicular magnetic field.

The twofold degenerate ground space of the effective Hamiltonian in Eq. (26) is spanned by the logical code words in Eq. (4), which are obtained from Eq. (38) by considering $\mathbf{k} = (0, 0)^T$ and $\mathbf{k} = (0, \sqrt{\pi})^T$, respectively. As discussed in Sec. IV B and in Appendix E, $\pi/2$ rotations in the $x_1 x_2$ plane correspond to a Fourier transform after the

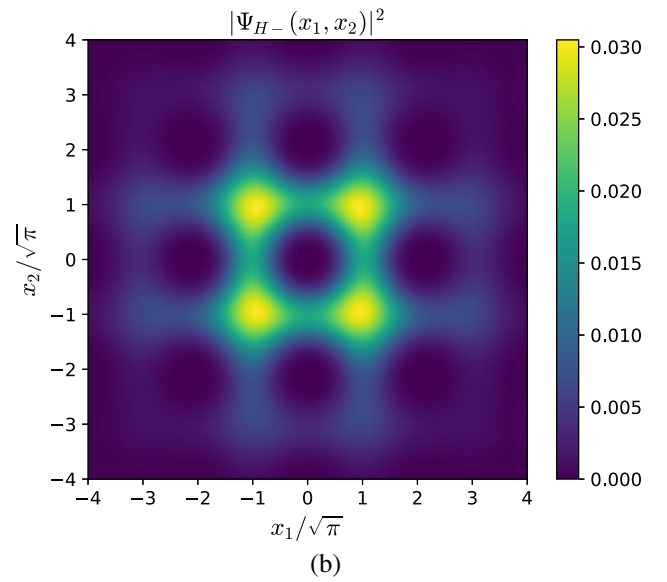
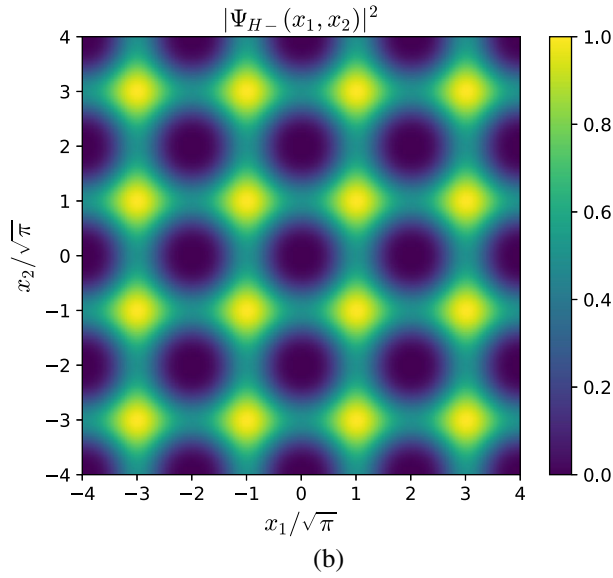
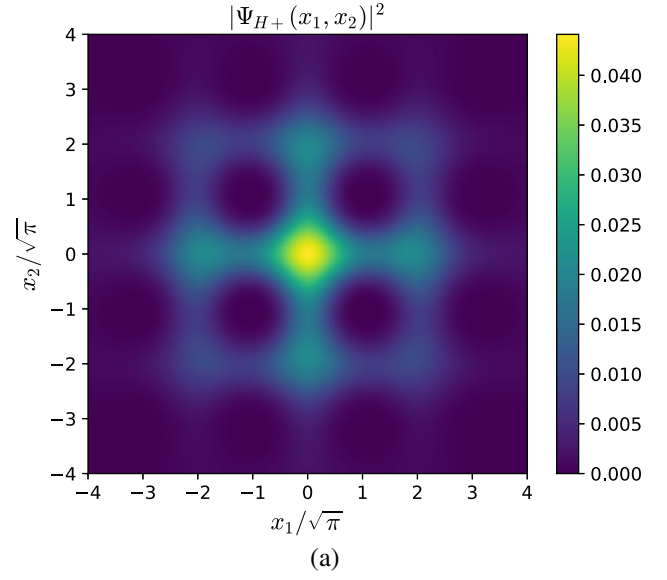
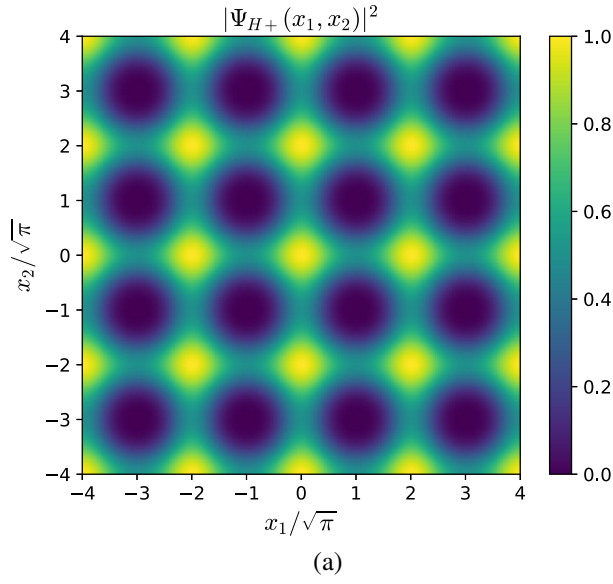


FIG. 8. Absolute values squared of the wave functions (a) $\Psi_{H+}(x_1, x_2)$ and (b) $\Psi_{H-}(x_1, x_2)$. These wave functions approximate the low-energy eigenstates of the two-dimensional Hamiltonian in Eq. (21) and are constructed to be even and odd, respectively, under a $\pi/2$ rotation in the x_1x_2 plane. Note that the wave functions in the plot are not normalized. We remark that $|\Psi_{H\pm}(x_1, x_2)|^2$ are the Husimi Q quasiprobability functions associated with the eigenstates $\psi_{H\pm}(X)$ [defined in Eq. (41)] of the projected Hamiltonian in Eq. (26).

LLL projection. Consequently, because in the basis $\psi_0(X)$ and $\psi_1(X)$ a Fourier transform is equivalent to a Hadamard gate [16] (see Sec. IV B), to construct fourfold rotational symmetric wave functions in the two-dimensional plane, we consider the linear combinations [91]

FIG. 9. Absolute values squared of the quasidegenerate ground state wave functions (a) $\Psi_{H+}(x_1, x_2)$ and (b) $\Psi_{H-}(x_1, x_2)$ of the Hamiltonian in Eq. (28), which includes a parabolic confinement potential. These functions are obtained by combining Eqs. (31) and (43) and are even and odd, respectively, under a $\pi/2$ rotation in the x_1x_2 plane. The wave functions here are normalized and are obtained by using $\Delta = 0.25$.

$$\psi_{H+}(X) = \cos\left(\frac{\pi}{8}\right)\psi_0(X) + \sin\left(\frac{\pi}{8}\right)\psi_1(X), \quad (41a)$$

$$\psi_{H-}(X) = -\sin\left(\frac{\pi}{8}\right)\psi_0(X) + \cos\left(\frac{\pi}{8}\right)\psi_1(X). \quad (41b)$$

These functions are even (odd) under Fourier transform, and so the corresponding two-dimensional wave functions

$\Psi_{H\pm}(x_1, x_2)$ are even (odd) under a $\pi/2$ rotation around the origin $x_i = 0$. The absolute values of the functions $\Psi_{H\pm}(x_1, x_2)$ are shown in Fig. 8.

We observe that the absolute values are periodic with period $2\sqrt{\pi}$ in both the x_1 and x_2 direction. Also, we find that these states are related to each other by

$$\Psi_{H+}(x_1, x_2) = ie^{i\sqrt{\pi}(x_2 - x_1)/2} \Psi_{H-}(x_1 + \sqrt{\pi}, x_2 + \sqrt{\pi}), \quad (42)$$

and so the absolute values of the two wave functions are simply obtained by a shift of $\sqrt{\pi}$ in the x_1 and x_2 direction.

As long as the Landau-level coupling remains weak, Eq. (36) is appropriate to describe also the system discussed in Sec. IV, where an additional parabolic potential is included. In particular, we find that the approximate grid states given in Eq. (32) transform into

$$\begin{aligned} \Psi_j(x_1, x_2) &= \sqrt{\frac{2\Delta^2}{\pi(1 + \Delta^2 + \Delta^4)}} e^{[\Delta^2(x_1 - ix_2)^2]/[2(1 + \Delta^2 + \Delta^4)]} e^{-[x_1(x_1 - ix_2)]/2} \\ &\times \vartheta\left[\frac{j}{2}\right]\left(\frac{-i(x_1 - ix_2)}{\sqrt{\pi(1 + \Delta^2 + \Delta^4)}}, 2i\frac{1 + \Delta^2}{1 + \Delta^2 + \Delta^4}\right), \end{aligned} \quad (43)$$

with $j = 0, 1$ and Δ being defined in Eq. (33). The low-energy eigenstates $\Psi_{H\pm}(x_1, x_2)$ of the Hamiltonian in Eq. (28) are related to these basis states by Eq. (31), and their absolute values obtained for $\Delta = 0.25$ are shown in Fig. 9.

Comparing Figs. 8 and 9, we observe that the parabolic potential introduces a Gaussian decay roughly of the order $1/\Delta^2$ of the wave functions in both the x_1 and x_2 direction and also it distorts the arguments of the theta functions with corrections of the order of Δ^2 . Of course, the absolute values of the wave functions in Fig. 8 are recovered by taking the limit $\Delta \rightarrow 0$.

VI. GKP HAMILTONIAN IN A NONRECIPROCAL SUPERCONDUCTING CIRCUIT

Here, we propose a possible experimental realization of the Hamiltonian in Eq. (28) based on a combination of state-of-the-art nonreciprocal superconducting circuits. We consider here the circuit shown in Fig. 10.

The device consists of two fluxonia coupled by a gyrator. The fluxonium is a well-known superconducting circuit comprising a Josephson junction with Josephson energy E_J in parallel with a capacitance C and an inductance L [92–94]. The crucial difference of our design from more conventional superconducting qubit architectures is the nonreciprocity that comes from the gyrator [39].

A gyrator is a two-port linear device that relates incoming currents and voltages according to

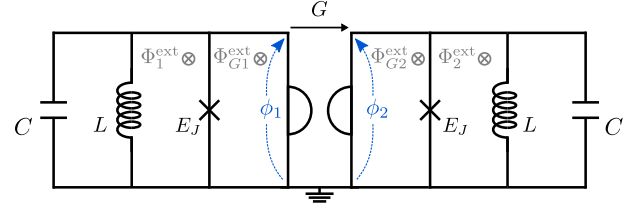


FIG. 10. Circuit design implementing the Hamiltonian in Eq. (46), which approximates the GKP Hamiltonian.

$$\underbrace{\begin{pmatrix} I_1 \\ I_2 \end{pmatrix}}_Y = \underbrace{\begin{pmatrix} 0 & -G \\ G & 0 \end{pmatrix}}_Y \begin{pmatrix} V_1 \\ V_2 \end{pmatrix}, \quad (44)$$

where G is the frequency-independent gyration conductance. Because it is characterized by an antisymmetric admittance matrix Y , this device is nonreciprocal and breaks the time-reversal symmetry of the circuit.

While the typical implementations of these devices are quite bulky [40,41], there are also recent realizations of miniaturized on-chip nonreciprocal devices based on actively pumped systems [42–45] or based on the quantum (anomalous) Hall effect [50,52]. Although our model is independent of the specific realization of the gyrator, the latter devices are advantageous in this context, because they are passive and they rely on quantized excitations with a long lifetime that can be well described by the theory of circuit quantum electrodynamics (cQED) [95,96]. A further advantage is the value of G .

To describe the system with circuit quantization theory, we introduce the node fluxes $\phi_i(t) = \int_{-\infty}^t V_i(t') dt'$. Using Kirchoff's laws, the following contribution to the Lagrangian [97,98]:

$$\mathcal{L}_G = \frac{G}{2} (\phi_1 \dot{\phi}_2 - \dot{\phi}_1 \phi_2) \quad (45)$$

correctly reproduces the defining property of the gyrator in Eq. (44) when two general classical networks are attached to it. Importantly, note that (in the style of an electronic system) Eq. (45) is similar to the effect of a homogeneous magnetic field of strength $B = G/e$ passing through the $\phi_1 \phi_2$ plane. More details on circuit quantization of nonreciprocal devices can be found in Refs. [97,99].

For now, we neglect the effect of magnetic fluxes threading the superconducting loops, and we set $\Phi_i^{\text{ext}} = \Phi_{Gi}^{\text{ext}} = 0$. Combining conventional circuit QED with Eq. (45), we find the Hamiltonian of the circuit in Fig. 10 to be

$$\begin{aligned} H &= \frac{(Q_1 + G\phi_2/2)^2}{2C} + \frac{(Q_2 - G\phi_1/2)^2}{2C} + \frac{1}{2L} (\phi_1^2 + \phi_2^2) \\ &- E_J \left[\cos\left(\frac{2\pi}{\Phi_{0,s}} \phi_1\right) + \cos\left(\frac{2\pi}{\Phi_{0,s}} \phi_2\right) \right]. \end{aligned} \quad (46)$$

TABLE I. Mapping of the parameters and variables used in the jargon of a crystal electron and cQED, such that the Hamiltonians in Eqs. (28) and (46) coincide. Note that the cyclotron frequency in terms of circuit parameters is $\omega_c = G/C$.

Crystal electron	x_i	p_i	m	eB	V	L_0	ω_0
Circuit	ϕ_i	Q_i	C	G	E_J	$\Phi_{0,s}$	ω_{LC}

We then impose the canonical commutation relation $[\phi_i, Q_j] = i\hbar\delta_{ij}$. Here, Q_i are the charges on the i th capacitor. Note that the superconducting flux quantum $\Phi_{0,s} = h/2e$ differs from the flux quantum Φ_0 used in the previous sections by a factor of 2. For simplicity, we assumed here that the two fluxonia coupled to the gyrator are identical. We do not expect small anisotropies to drastically alter the results described in this section.

The Hamiltonian in Eq. (46) describing our circuit has the same structure as the Hamiltonian of a confined crystal electron in a magnetic field in Eq. (28) and discussed in detail in Sec. IV. The variables that play equivalent roles in the two cases are given in Table I.

In particular, the gyration conductance G acts as a magnetic field B , and the characteristic frequency of the LC circuit $\omega_{LC} = 1/\sqrt{LC}$ acts as the harmonic confinement ω_0 . For later convenience, we also introduce the charging energy $E_C = e^2/2C$ and the inductive energy $E_L = \Phi_{0,s}^2/4\pi^2L$.

As shown in Sec. III, the number of flux quanta threading one unit cell is of fundamental importance for realizing GKP states. In our circuit, Eq. (17) becomes

$$\frac{p}{q} = \frac{G}{e} \frac{\Phi_{0,s}^2}{\Phi_0} = \frac{G}{G_0}, \quad (47)$$

where we introduce the superconducting conductance quantum $G_0 = (2e)^2/h$. In order to obtain GKP states, we require $p/q = 1/2$, and, accordingly, we require the gyration conductance to be precisely

$$G = \frac{2e^2}{h}. \quad (48)$$

We remark again that, while this value of $G \sim 1/(13 \text{ k}\Omega)$ is unrealistic for superconducting-based gyrators, it can be easily reached using quantum (anomalous) Hall effect devices, where the characteristic impedance is $1/G = h/2e^2\nu$ [46–52], with ν being the Landau-level filling factor. The robust quantization of the Hall conductivity in these materials also guarantees that the value of G remains precisely fixed for a wide range of design parameters, hence improving the reproducibility of the gyrator.

To reach low values of the harmonic confinement frequency ω_{LC} , we expect that the novel hyperinductances [100], the kinetic inductances based on granular aluminum

TABLE II. Design parameters for the circuit in Fig. 10. These parameters are achievable in state-of-the-art superconducting circuits. The value of the inductive energy is taken from Ref. [100]. The charging energy and the inductive energy correspond to a capacitance $C = 1.4 \text{ fF}$ and an inductance $L = 2.3 \text{ }\mu\text{H}$, respectively.

Parameter	GHz
E_C/h	13.50
E_J/h	3.50
E_L/h	0.07
$\omega_c/2\pi$	8.59
$\omega_{LC}/2\pi$	2.75

[101,102], or thin Nb nanowires [103] will be suitable. Also, the Josephson junctions should work in the charge regime $E_C \gtrsim E_J$, which guarantees a weak Landau-level coupling. In Table II, we list parameter values that are experimentally achievable in state-of-the-art superconducting circuits and that can be used to design GKP qubits. The resulting, relevant energy ratios which need to be small are $E_J/\hbar\omega_c = 0.4$ and $\hbar\omega_{LC}/E_J = 0.8$. For the parameter defining the widths of the approximate grid states (see Sec. IV B), we obtain $\Delta = (E_L/E_J)^{1/4}e^{\pi/4} = 0.8$. In Sec. VII, we discuss how these parameters should be improved and also which trade-offs might be needed to operate the circuit in the desired way.

We emphasize that our circuit encodes the approximate grid states in a subsystem (to be precise, in the LLL) whose dynamics is effectively described by the approximate GKP Hamiltonian in Eq. (30). For this reason, the code words are passively protected [22], and so, in contrast to current efforts to encode grid states in superconducting cavities [18], they do not require permanent active stabilization [22].

We also point out that there is a different proposal for a superconducting circuit implementing grid states in a doubly nonlinear qubit (the dualmon) [21], which involves a Josephson junction and a quantum phase slip wire. However, in contrast to our proposal, its dynamics is not described within a Landau-level projection, and also the GKP code words are not the lowest-lying eigenstates of the resulting Hamiltonian [21].

So far, we neglect the effect of potential external magnetic fluxes threading the superconducting loops in the circuit shown in Fig. 10. Also, no external voltage or current sources are attached to it. In the following, we show how these additional degrees of freedom can be used to perform single- and two-qubit gates and for state preparation and qubit readout. Because of the equivalent roles of X and P , stemming from the similar accessibility of both the fluxes ϕ_1 and ϕ_2 in the circuit in Fig. 11, quantum operations can be performed in the logical $\bar{Z}$ or $\bar{X}$ basis, depending on to which port of the gyrator the sources are applied.

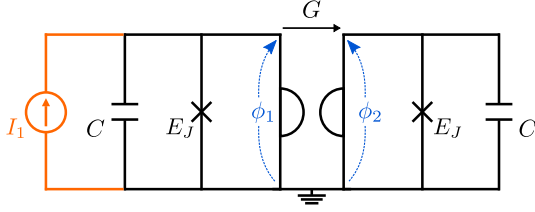


FIG. 11. Circuit implementation of the *ideal* logical $\bar{Z}$ gate with a current source. Assuming the system starts in the code subspace, the current source is turned on and kept constant at a value I_1 for a time t_Z given in Eq. (54). A current source on the opposite port of the gyrator would instead implement a logical $\bar{X}$ gate.

A. Logical $\bar{X}$ and $\bar{Z}$ gates

We now turn our attention to the implementation of logical gates in our system by focusing on the single-qubit $\bar{X}$ and $\bar{Z}$ gates defined in Eq. (3). For the sake of clarity, we carry out the analysis for the case without harmonic confinement potential, i.e., without inductances in the superconducting circuit. The same procedures also work for the complete circuit in Fig. 10 when the ratios E_L/E_J and E_J/E_C are sufficiently small, yielding approximate logical gates.

In our analysis, we demand that the system is operated in the relevant case of weak Landau-level coupling as discussed in Sec. III A. The logical operators $\bar{X}$ and $\bar{Z}$ can be implemented using current sources shunting either of the ports of the gyrator. The circuit implementing the $\bar{Z}$ gate is depicted in Fig. 11.

The Hamiltonian of this circuit can be written as

$$H(t) = \frac{\pi^2}{2C} - E_J \cos \left[\frac{2\pi}{\Phi_{0,s}} \left(R_1 + \frac{\pi_2}{C\omega_c} \right) \right] - E_J \cos \left[\frac{2\pi}{\Phi_{0,s}} \left(R_2 - \frac{\pi_1}{C\omega_c} \right) \right] - I_1(t) \left(R_1 + \frac{\pi_2}{C\omega_c} \right), \quad (49)$$

where, in analogy with the electronic case [see Eqs. (11) and (12)], we define the dynamical momenta $\pi_1 = Q_1 + G\phi_2/2$ and $\pi_2 = Q_2 - G\phi_1/2$ and the guiding centers $R_1 = \phi_1 - \pi_2/C\omega_c$ and $R_2 = \phi_2 + \pi_1/C\omega_c$. The current source appears in the Hamiltonian in the term $-I_1(t)[R_1 + \pi_2/C\omega_c] = -I_1(t)\phi_1$. Note that, in the electronic analogy, the current source acts as a homogeneous and time-dependent in-plane electric field, whose direction depends on the port the generator is connected to. We expect that, if the current source $I_1(t)$ does not contain frequencies close to ω_c , it will not cause transitions between Landau levels. Hence, we can project onto the LLL and obtain (dropping constant shifts in energy)

$$H_{\text{LLL}} = -V_0[\cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)] - \frac{I_1(t)\Phi_{0,s}}{\sqrt{\pi}}X = H_{\text{GKP}} - \frac{I_1(t)\Phi_{0,s}}{\sqrt{\pi}}X, \quad (50)$$

where we immediately perform the variable rescaling in Eq. (25). Also, we use the definition of the GKP Hamiltonian in Eq. (7) to identify $V_0 = E_J e^{-\pi}$.

Now, we consider the following scenario. At time $t = 0$, the state is assumed to be in a generic superposition $|\psi_{\text{in}}\rangle = c_0|\bar{0}\rangle + c_1|\bar{1}\rangle$ of the ideal GKP code words given in Eq. (4). We assume a constant current source $I_1(t) \equiv I_1$. In this case, the time evolution operator associated with H_{LLL} in Eq. (50) reduces to $U_{\text{LLL}}(t) = \exp(-iH_{\text{LLL}}t/\hbar)$. In order to understand the effect of $U_{\text{LLL}}(t)$ on $|\psi_{\text{in}}\rangle$, we use the Zassenhaus formula [104]

$$e^{t(A+B)} = e^{tA} e^{tB} e^{-t^2[A,B]/2} e^{t^3(2[B,[A,B]] + [A,[A,B]])/6} \dots \quad (51)$$

with $A = iI_1\Phi_{0,s}X/\hbar\sqrt{\pi}$ and $B = -iH_{\text{GKP}}/\hbar$. Since all the commutators in the Zassenhaus formula have the GKP states as degenerate eigenstates, e.g.,

$$[X, \cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)] = -i2\sqrt{\pi} \sin(2\sqrt{\pi}P), \quad (52)$$

one can show that

$$|\psi(t)\rangle = U_{\text{LLL}}(t)|\psi_{\text{in}}\rangle = e^{i\theta(t)} \exp\left(i \frac{I_1\Phi_{0,s}}{\hbar\sqrt{\pi}} X t\right) |\psi_{\text{in}}\rangle, \quad (53)$$

where we factorize the irrelevant phase factor $e^{i\theta(t)}$. Thus, after a time

$$t_Z = \frac{\hbar\pi}{I_1\Phi_{0,s}}, \quad (54)$$

a logical $\bar{Z}$ gate is applied to $|\psi_{\text{in}}\rangle$, up to an overall phase. Ideally, after a time t_Z , the current source must be switched off.

Because the variables X and P refer to the left and the right port of the gyrator, respectively, it is straightforward to convert between the logical $\bar{Z}$ and $\bar{X}$ basis by simply changing the port of the gyrator where the current source is applied and using the same protocol.

B. Noise sensitivity

We provide a first analysis of the noise sensitivity of our qubit to typical noise sources, such as flux and charge noise. We start our discussion by analyzing charge noise. In the circuit in Fig. 10, charge noise can be modeled by capacitively coupling random voltage sources to the ports of the gyrator. This modifications change the kinetic term in the Hamiltonian in Eq. (46) as $\pi^2/2C \mapsto (\pi + \mathbf{Q}_g)^2/2C$, where $\mathbf{Q}_g$ is a vector containing the random charges on the capacitors connected to the voltage sources on either side of the circuit. The eigenspectrum is insensitive to static gate

charges, since they can be gauged away by a unitary transformation, as for the fluxonium qubit [92,105]. Moreover, charge noise couples to the dynamical momenta $\pi_{1,2}$, and, as a consequence, it has only a small effect on the guiding center variables in which our states are encoded. From these arguments, we conclude that charge noise should not be a major source of decoherence in our system, even if the transmon condition is not fulfilled [106].

Another typical noise source in our system is flux noise. We begin our analysis of flux noise sensitivity by considering again the ideal GKP Hamiltonian defined in Eq. (7), thus neglecting the confining potential of the inductive shunts. After the LLL projection, the external fluxes through the loops formed by gyrator branches and Josephson junctions give rise to the Hamiltonian [107]

$$H/V_0 = -\cos[2\sqrt{\pi}X + \varphi_{G1}^{\text{ext}}(t)] - \cos[2\sqrt{\pi}P + \varphi_{G2}^{\text{ext}}(t)], \quad (55)$$

where $\varphi_{G1,G2}^{\text{ext}}(t) = 2\pi\Phi_{G1,G2}^{\text{ext}}(t)/\Phi_{0,s}$ are the reduced magnetic fluxes through the loops on either port of the gyrator, respectively. The GKP code space is intrinsically protected with respect to these fluxes, independently of their strengths. In order to show this protection, we rewrite Eq. (55) as the sum of the desired GKP Hamiltonian and additional noise operators with time-dependent coefficients, i.e.,

$$H/V_0 = H_{\text{GKP}}/V_0 + s_X(t) \sin(2\sqrt{\pi}X) + c_X(t) \cos(2\sqrt{\pi}X) + s_P(t) \sin(2\sqrt{\pi}P) + c_P(t) \cos(2\sqrt{\pi}P), \quad (56)$$

where we define $c_{X,P}(t) = 1 - \cos[\varphi_{G1,G2}^{\text{ext}}(t)]$ and $s_{X,P}(t) = \sin[\varphi_{G1,G2}^{\text{ext}}(t)]$. Because all the individual noise operators in Eq. (56) have the GKP code space as degenerate eigensubspace, we conclude that the GKP code space is a decoherence-free subspace [108,109] with respect to this kind of noise. A similar observation is also made for the dualmon in Ref. [21]. We stress that this argument does not rely on the assumption of Markovianity of the flux noise.

The previous derivation assumes the ideal GKP Hamiltonian given in Eq. (7). However, when we work with its confined version, corresponding to the circuit shown in Fig. 10, we have to take into account the noise associated with the external magnetic fluxes $\Phi_{1,2}^{\text{ext}}$ through the superconducting loops formed by the Josephson junctions and the inductances on each port of the gyrator. We stress that, in the limit of large inductances that we are considering here, this flux noise has a weak effect and that the associated noise term vanishes as the value of the inductances increases.

To avoid pure dephasing due to $1/f$ noise, such as flux noise, the energy levels should not depend on the noise parameters. An example of the dependence of the energy levels on flux noise is shown in Fig. 12. The protection

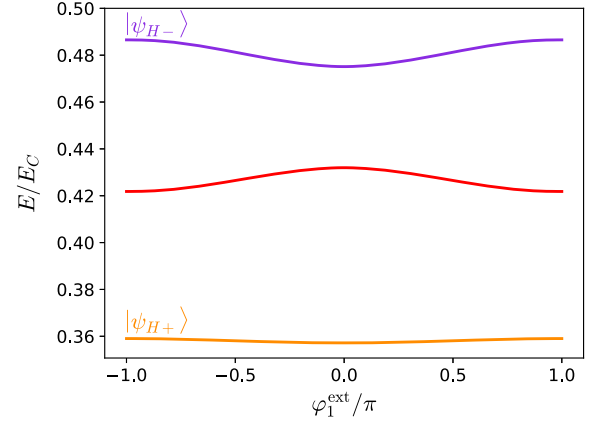


FIG. 12. Low-energy spectrum of the circuit shown in Fig. 10 as a function of the external flux $\varphi_1^{\text{ext}} = 2\pi\Phi_1^{\text{ext}}/\Phi_{0,s}$ for the parameters in Table II and with the other external fluxes set to zero. For these values, $|\psi_{H-}\rangle$ is the second excited state. One can clearly notice the sweet spots when φ_1^{ext} is an integer multiple of π .

against flux noise dephasing, for the realistic parameter regime in Table II, does not seem to be inherently different from that of state-of-the-art fluxonium qubits [92], as well as the one of $0-\pi$ qubits [58] and bifluxon qubits [110], where the energy levels show a behavior as a function of the external fluxes similar to our qubit.

However, in analogy to the $0-\pi$ qubit [58] and the bifluxon [110], the disjoint support of the GKP code words with respect to both X and P guarantees that the matrix elements of local noise operators between the encoded states are very small. This smallness, in turn, guarantees protection against energy relaxation. In our case, the relevant operators to characterize the noise due to the external fluxes are $\varphi_i = 2\pi\phi_i/\Phi_{0,s}$ (coupling to Φ_i^{ext} and Φ_{Gi}^{ext}) and, to first order in the small noise parameters, $\sin(\varphi_i)$ (coupling to Φ_{Gi}^{ext}). Furthermore, the noise operator associated with quasiparticle tunneling is $\sin(\varphi_i/2)$ [111]. These noise sources have a small effect in the relevant parameter regime. In fact, because the wave functions $\psi_{H\pm}(X)$ are both even in X , the matrix elements of the noise operators in the LLL projection vanish in the subspace spanned by $\psi_{H\pm}(X)$, i.e.,

$$\langle i | \Pi_{\text{LLL}} \mathcal{O}_{\text{noise}} \Pi_{\text{LLL}} | j \rangle = 0, \quad (57)$$

with $\mathcal{O}_{\text{noise}} \in \{\varphi_{1,2}, \sin(\varphi_{1,2}), \sin(\varphi_{1,2}/2)\}$ and $i, j \in \{\psi_{H+}, \psi_{H-}\}$.

Note, however, that for the realistic, currently achievable parameter regime in Table II, there is a level crossing of the first and second excited states as the inductances decrease; see also Fig. 4. Although this crossing also happens for the $0-\pi$ qubit [60] for the corresponding experimentally realized parameters, we believe that this undesired level crossing could be more problematic in our system. In fact,

the ground state and the second excited state both couple to this intermediate level (for instance, through the flux variable), which could lead to a decay of the second excited state into the ground state via this level. However, considering a value of E_L that is 5 times smaller than the one listed in Table II, the undesired energy level is pushed above the energies of the code words, thus solving this issue.

C. State preparation, qubit readout, and Clifford gates

State preparation, qubit readout, and the implementation of Clifford gates are nontrivial and related topics for our qubit. A destructive measurement in the GKP basis $\{|\bar{0}\rangle, |\bar{1}\rangle\}$ can be performed by measuring the flux ϕ_1 , which in the LLL projection is approximately equivalent to measuring the rescaled guiding center variable X . An outcome of the measurement that is close to an even multiple of $\sqrt{\pi}$ corresponds to state $|\bar{0}\rangle$, while an outcome close to an odd multiple of $\sqrt{\pi}$ is assigned to state $|\bar{1}\rangle$.

A nondestructive measurement can be instead implemented if we have the ability to perform a GKP phase estimation protocol [112], where we prepare an ancilla qubit in $|\bar{0}\rangle$, perform a CNOT with the ancilla qubit as target, and then measure the ancilla destructively [113,114]. In this protocol, the nondestructive measurement relies on the possibility to prepare the logical state $|\bar{0}\rangle$.

As recently shown in Ref. [54], a logical $|\bar{0}\rangle$ state can be prepared deterministically by an adaptive protocol starting from two Hadamard eigenstates, using Clifford operations and a destructive readout of one of the two qubits. In Sec. IV B, we show that in the symmetric case the ground state of our system is indeed approximately a GKP Hadamard eigenstate. Thus, the ability to cool down our system in the ground state would give us also the ability to prepare the GKP $|\bar{0}\rangle$ state, when combined with Clifford operations and destructive measurement described above. We also remark that, in Ref. [85], it is shown instead how to prepare the GKP Hadamard eigenstate starting from many GKP logical $|\bar{0}\rangle$ states. However, as we observe in Sec. IV B, if we have the ability to tune the asymmetry between the fluxonia, for instance, by considering sufficiently asymmetric Josephson junctions, the ground state becomes an approximate GKP $|\bar{0}\rangle$ state. Hence, if we are able to prepare this state, this preparation would be an alternative to using the protocols described beforehand.

In Sec. VI A, we describe a protocol to implement logical $\bar{X}$ and $\bar{Z}$ gates by means of current sources. Here, we discuss further ideas for the implementation of general Clifford operations. As discussed in Ref. [16], one of the convenient properties of the GKP code is that, in the encoded subspace, Clifford unitaries are implemented by symplectic transformations (see also Ref. [115] for a review of gates for the GKP code). Symplectic transformations are generated by Gaussian unitaries and, as such, can be

realized by using linear optics and squeezing [55,116]. The Clifford group for a single qubit is generated by the Hadamard gate $\bar{H}$ defined in Eq. (34) and the phase gate [117]

$$\bar{S} = \begin{pmatrix} 1 & 0 \\ 0 & i \end{pmatrix}. \quad (58)$$

As discussed in Ref. [22], a possible way to implement the phase gate in our GKP qubits relies on the ability to change the magnitude of one of the quadratic terms in Eq. (30). This change can be achieved by tuning in time the superinductances, effectively creating an asymmetry between the two fluxonia in the circuit in Fig. 10. In fact, in the GKP code, the ideal unitary implementing the phase gate can be chosen as $U_{\bar{S}} = e^{-iX^2/2}$ [16,115], and so we need a term $\propto X^2$ that dominates the quadratic part of the Hamiltonian. This term appears if we create an asymmetry between the two inductances such that $E_{L_1} \gg E_{L_2}$. Then, in the LLL projection, we obtain a quadratic term $\propto E_{L_1}X^2$ that dominates over $\propto E_{L_2}P^2$.

A logical Hadamard gate $\bar{H}$ can be implemented with a similar protocol by changing the magnitude of both the quadratic terms in Eq. (30), e.g., by varying both the superinductances in the circuit in Fig. 10 in time. However, the Hadamard gate $\bar{H}$ in Eq. (34), that corresponds to a change of basis from logical $\bar{Z}$ to logical $\bar{X}$, is approximately equal to a Fourier transform. Thus, the Hadamard gate $\bar{H}$ can be also realized virtually, namely, by applying consequent gates and measurements on the other port of the gyrator. This procedure is used in Sec. VI A for the discussion of the $\bar{Z}$ and $\bar{X}$ gates, since $\bar{X} = \bar{H}\bar{Z}\bar{H}$.

Similar ideas can also be employed to implement a CNOT gate: In this case, we need a tunable inductance coupling the branches of two of our GKP qubits. In addition, for the experimental realizable parameters in Table II, we believe that, in analogy to the $0-\pi$ qubit, one could perform gates also by using higher excitations of the circuit [59]. These ideas have been recently realized experimentally for the $0-\pi$ qubit [60].

The preparation of the ground state becomes more and more difficult as the quality of our GKP states improves. In fact, as the inductances increase, the ground and first excited states become closer in energy. In this case, state preparation would require temperatures that are lower than in current practice for superconducting qubits. Similar problems emerge also in the heavy fluxonium qubit and have been recently addressed in Ref. [118]. Akin to the implementation of Clifford gates, an alternative approach to state preparation could use tunable superinductances. In this scheme, one prepares the ground state at relatively small inductances and then adiabatically increases the inductances keeping the system always in the ground state; see Fig. 4. However, also this scheme becomes harder as

the energy gap shrinks, but the fact that the two quasidegenerate states are not significantly coupled by terms in the Hamiltonian should suppress transitions from one state to the other in the adiabatic protocol.

We believe that similar protocols can be realized by using superconducting quantum interference devices (SQUIDs) [119,120] instead of Josephson junctions and by modifying the effective Josephson energy by adiabatically tuning the external fluxes. Here, we do not explore further these protocols quantitatively and leave a detailed description of the implementation to future research.

VII. COMPARISON WITH OTHER SUPERCONDUCTING QUBITS AND DESIGN TRADE-OFFS

Our system ranks in the class of superconducting qubits that attempt to provide a built-in, passive protection against noise, going beyond standard qubit designs such as the transmon [106], the flux qubit [121], and the fluxonium [92]. The prototypical example of a protected superconducting qubit is perhaps the $0-\pi$ qubit, originally proposed in Ref. [20] (with ideas from Ref. [19]), further analyzed in Refs. [57–59], and recently also experimentally realized [60]. Reference [61] instead proposes a different circuit design to realize essentially the same idea of the original $0-\pi$ qubit circuit. The $0-\pi$ qubit achieves the desired protection by engineering a superconducting circuit whose energy levels are robust against typical noise sources and whose eigenstates have disjoint support, thus inhibiting dissipation-induced transitions. The same idea of protection is employed also in the recently proposed, and experimentally realized, bifluxon [110].

Among these qubit designs, the system that shows more similarities with our proposal, although realized by a completely different superconducting circuit, is certainly the proposed dualmon [21]. The dualmon is constructed by shunting a Josephson junction in parallel to a quantum phase slip wire, and its effective dynamics is described by the Hamiltonian in Eq. (26), however, for $p/q = 1$ [122]. Thus, the crucial difference to our proposal is that the ground state of the dualmon is unique and nondegenerate. Nonetheless, the dualmon still possesses states that are robust against flux and charge noise, in which quantum information can be reliably encoded. Also, we remark that in our case the quantum information is encoded in a subsystem, while the other degrees of freedom are assumed to be frozen in a Landau level.

The experimental realization of our system relies on the availability of on-chip gyrators with gyration conductance being twice the conductance quantum, i.e., $G = 2e^2/h$. This realization will be most likely achievable by using quantum Hall devices [46–52]. Assuming that this realization is possible, we reported in Table II currently achievable parameters in superconducting systems, for which the Landau-level approximation is valid.

We remark that, in order to fully exploit the potentials of the proposed circuit design in Fig. 10, an improvement of the currently achievable parameters is needed. In particular, a decrease of the ratio E_L/E_J by 2 orders of magnitude (while keeping the Landau-level projection valid) seems necessary in order to obtain adequate approximate grid states with a small width Δ . We observe that reaching sufficiently small values of Δ is partially hindered by the unfavorable rescaling of the parameters in the lowest Landau-level projection. Note also that a considerable decrease of the ratio E_L/E_J cannot be simply achieved by increasing the Josephson energy E_J , since (given the fact that G is fixed) the main parameter that controls the validity of the Landau-level projection is E_J/E_C . Thus, the requirement of a smaller value of E_L/E_J goes in the direction of increasing the superinductance, which is a desirable feature also for the previously discussed protected qubits.

In this work, we show the connection with the GKP Hamiltonian, and, thus, we focus on the parameter regime in which the Landau-level projection holds. However, we believe that the system could also show some different forms of noise resilience when this approximation is not valid. The study of the system in a different parameter regime, especially in an experimentally easily realizable one, is definitely an interesting topic for future work.

VIII. CONCLUSIONS AND OUTLOOK

We have designed a circuit composed of state-of-the-art superconducting circuit elements and a nonreciprocal device, that can be used to passively implement the GKP quantum error-correcting code. Our proposal crucially relies on the gyrator, which plays the role of an effective homogeneous magnetic field in an analogous electronic system and whose amplitude depends on the characteristic admittance of the device. By taking advantage of recent advances in manufacturing nonreciprocal quantum Hall effect devices, one can reliably reach the desired value of the effective magnetic field, which is well outside of the range that can be obtained in electronic systems.

By working out in detail the equivalence between our circuit and the problem of an electron in a magnetic field in a crystal potential, we analyze the system and identify a parameter range where the ground states of the system are the GKP code words. Our analysis shows the deep relation between the GKP states and the Hofstadter butterfly, which, to the best of our knowledge, was not known previously. We study an implementation of approximate GKP code words by shunting our circuit with large inductances.

We work out a mapping that allows one to understand the eigenstates of the system in different coordinate systems, facilitating the interpretation of experimental results.

We discuss possible ways to implement one- and two-qubit logical gates as well as ideas for state preparation and qubit readout. This discussion suggests that universal quantum computation can be done with our qubits by using

only current sources and tunable inductances or tunable Josephson junctions (SQUIDS).

Finally, we discuss the effect of typical noise sources, i.e., charge and flux noise, and conclude that our qubit is well protected against them.

In this paper, we list a few ideas of how to implement phase gates and how to initialize the quantum state. A detailed comparison between the different protocols is still missing and is required to have a better understanding of the experimental capability of our qubit. Also, a more realistic modeling of the device would have to account for asymmetries in the circuit, for a finite bandwidth of the gyrator or its internal degrees of freedom, whose effects have been overlooked in our analysis. A first consideration of a finite but appropriately large bandwidth indicates that the effective dynamics of the resulting system qualitatively reduces to the model described in this work. Thus, we believe that these imperfections in the experiments would not affect the qualitative behavior of the system, which provides a promising hardware implementation of the GKP code.

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APPENDIX A: MAGNETIC TRANSLATION OPERATORS

In this Appendix, we summarize a few key results about the magnetic translation operators (MTOs), which are required in Sec. III. In analogy to the main text, here we restrict ourselves to the analysis of rational magnetic fluxes $\Phi/\Phi_0 = p/q$ [see Eq. (17)], where $\Phi = BL_0^2$ denotes the flux threading one unit cell of size $L_0 \times L_0$ and $\Phi_0 = h/e$ is the magnetic flux quantum. Using the definition of the MTOs in Eq. (14), for integer values of p and q we find

$$[T_1(qL_0), T_2(L_0)] = 0, \quad (\text{A1})$$

since the magnetic unit cell of size $qL_0 \times L_0$ contains p flux quanta. Equation (A1) justifies the magnetic Bloch theorem in Eq. (18), which defines the magnetic Bloch states $|\mathbf{k}\rangle$, with $\mathbf{k}$ restricted to the first Brillouin zone; see Eq. (19).

Given the magnetic Bloch theorem in Eq. (18), we can find basis states that describe the system within a unit cell by considering the eigenstates of the smallest possible

translations compatible with the magnetic Bloch theorem. The choice of operators is, of course, nonunique, and here, for example, we choose the eigenstate of the operator $T_2(L_0/p)$, which commutes with both $T_1(qL_0)$ and $T_2(L_0)$. We then define the basis states

$$T_2(L_0/p)|\mathbf{k}, l\rangle = e^{i(k_y L_0 + 2\pi l)/p} |\mathbf{k}, l\rangle, \quad (\text{A2})$$

where $l = 0, 1, 2, \dots, p-1$.

Note that the Hamiltonian in Eq. (21) exclusively comprises the MTOs $T_1(qL_0/p)$ and $T_2(qL_0/p)$ and their Hermitian conjugate. The action of $T_2(qL_0/p) = [T_2(L_0/p)]^q$ on $|\mathbf{k}, l\rangle$ follows straightforwardly from Eq. (A2). Thus, it remains to analyze the action of $T_1(qL_0/p)$. From Eq. (16), we find

$$\begin{aligned} T_2(L_0/p)T_1(qL_0/p)|\mathbf{k}, l\rangle \\ = e^{i[k_2 L_0 + 2\pi(l+1)]/p} T_1(qL_0/p)|\mathbf{k}, l\rangle, \end{aligned} \quad (\text{A3})$$

from which we conclude that

$$T_1(qL_0/p)|\mathbf{k}, l\rangle = e^{ik_1 q L_0/p} |\mathbf{k}, (l+1) \bmod p\rangle. \quad (\text{A4})$$

Note that the state $|\mathbf{k}, l\rangle$ maps into itself after p consequent applications of $T_1(qL_0/p)$, in agreement with Eq. (18).

Finally, we analyze the degeneracy of the eigenstates of the Hamiltonian H in Eq. (21). To this end, we note that $T_1(L_0)$ commutes with H but not with both the MTOs of the boundary conditions in Eq. (18). Thus, if $|\psi\rangle$ is an eigenstate of H with eigenenergy E , the state $T_1(L_0)|\psi\rangle$ is also an eigenstate of H with the same eigenenergy, and, because $[T_1(L_0), T_2(L_0)] \neq 0$, the states $|\psi\rangle$ and $T_1(L_0)|\psi\rangle$ are physically distinguishable for $q > 1$. In particular, one can easily show that

$$\begin{aligned} T_1(L_0)|k_1, k_2\rangle^T, l\rangle \\ = e^{ik_1 L_0} |[k_1, (k_2 + 2\pi/q L_0) \bmod 2\pi/L_0]^T, l\rangle \\ \propto |(k_1, k_2)^T, l\rangle, \quad \text{for } q > 1. \end{aligned} \quad (\text{A5})$$

As a result, every energy band of the Hamiltonian is at least q -fold degenerate.

APPENDIX B: NUMERICAL ANALYSIS OF THE EIGENSYSTEM

1. Without confinement potential

In the following, we provide a method to numerically determine the spectrum of the Hamiltonian in Eq. (21). To this end, we expand the Hamiltonian in the product state basis $|n; \mathbf{k}, l\rangle$ [defined in Eqs. (22) and (23)] with $l = 0, \dots, p-1$ and $n = 0, \dots, N$ for some reasonably large integer N .

In particular, the matrix elements of the displacement operator in the Landau-level basis are known analytically [123] and read

$$\langle m|D_a(\alpha)|n\rangle = \sqrt{\frac{n!}{m!}}\alpha^{m-n}e^{-|\alpha|^2/2}L_n^{m-n}(|\alpha|^2), \quad (\text{B1})$$

where $L_n^{m-n}(|\alpha|^2)$ denotes the associated Laguerre polynomial. Note that the evaluation of the Hamiltonian in the given basis is particularly convenient in the weak Landau-level coupling limit ($V/\hbar\omega_c \ll 1$), since the coupling of product states $|n; \mathbf{k}, l\rangle$ with different n is weak. In this limit, every Landau level splits into p bands with finite widths, which are well separated from all the other split Landau levels; see Fig. 2. Considering only one Landau level, it is worth mentioning that the way in which it splits results in an energy spectrum which shows a fractal behavior similar to a deformed Hofstadter butterfly [34]; see Fig. 3.

2. Including confinement potential

Here, we present a convenient basis for the numerical analysis of the eigensystem of the Hamiltonian in Eq. (28). To this end, we introduce the bosonic ladder operators associated with the guiding center variables:

$$b = \frac{1}{\sqrt{2}} \frac{R_1 + iR_2}{l_B}, \quad b^\dagger = \frac{1}{\sqrt{2}} \frac{R_1 - iR_2}{l_B}, \quad (\text{B2})$$

satisfying $[b, b^\dagger] = 1$, and define the unitary displacement operator associated to these variables:

$$D_b(\beta) = e^{\beta b^\dagger - \beta^* b}, \quad \beta \in \mathbb{C}. \quad (\text{B3})$$

Given the ladder operators of the guiding center variables and those of the dynamical momenta [see Eq. (20)], we rewrite the Hamiltonian in Eq. (28) as (dropping constant energy offsets)

$$H = \hbar\omega_c a^\dagger a + \frac{\hbar\omega_0^2}{\omega_c} (a^\dagger a + b^\dagger b + ab + a^\dagger b^\dagger) - \frac{V}{2} [D_a(\lambda)D_b(-\lambda) + D_a(i\lambda)D_b(i\lambda) + \text{H.c.}], \quad (\text{B4})$$

with $\lambda = \sqrt{q\pi/p}$ being the absolute value of each displacement. This Hamiltonian is expanded in the basis of the Fock product states

$$|n, m\rangle = \frac{a^{\dagger n}}{\sqrt{n!}}|0\rangle \otimes \frac{b^{\dagger m}}{\sqrt{m!}}|0\rangle, \quad n, m \in \mathbb{N}_0, \quad (\text{B5})$$

where we have to reasonably truncate n and m . In the process, the matrix elements of each individual term in the Hamiltonian are known analytically, especially the matrix elements of the displacement operators; see Eq. (B1).

At this point, one could proceed with an analytical diagonalization of the quadratic part [124] of the Hamiltonian [first line in Eq. (B4)] in order to reduce the coupling of the basis states. The eigenstates of the

quadratic part of the Hamiltonian are known as Fock-Darwin states [86,87].

We, however, do not perform this diagonalization, because we want to retain the jargon of Landau levels. Nevertheless, both approaches coincide in the limit of consideration.

After the LLL projection, i.e., restricting to the subspace spanned by $|0, m\rangle$, the Hamiltonian in Eq. (B4) reduces to (dropping constant energy offsets)

$$H_{\text{LLL}} = \frac{\hbar\omega_0^2}{\omega_c} b^\dagger b - \frac{V_0}{2} [D_b(i\lambda) + D_b(-i\lambda) + D_b(\lambda) + D_b(-\lambda)], \quad (\text{B6})$$

where $V_0 = Ve^{-\pi q/2p}$. Note that, in the limit of weak confinements ($\hbar\omega_0^2/\omega_c V_0 \ll 1$), states $|0, m\rangle$ with different m are strongly coupled due to the crystal potential. Therefore, a large number of Fock states is required for an accurate numerical treatment. Nevertheless, a finite confinement prevents the eigenstates of constituting arbitrarily high excited Fock states.

Moreover, the matrix elements $\langle 0, m_1 | H_{\text{LLL}} | 0, m_2 \rangle$ are nonzero only if $(m_1 - m_2) \bmod 4 = 0$. Thus, the Hamiltonian in the LLL projection couples only every fourth Fock state. For this reason, also the eigenstates of H_{LLL} comprise only every fourth Fock state [16].

The consequence of this characteristic of the eigenstates becomes clear by considering the X representation [see Eq. (25)] of the m th Fock state:

$$\langle X | m \rangle = \frac{1}{\sqrt[4]{\pi}} \frac{1}{\sqrt{2^m m!}} H_m(X) e^{-X^2/2}, \quad (\text{B7})$$

where $H_m(X)$ is the m th Hermite polynomial. The Hermite functions in Eq. (B7) have the fundamental property of being eigenfunctions of the Fourier transform with eigenvalue $(-i)^m$ [125], which is cyclic in m with periodicity 4. Hence, we can conclude that also the eigenfunctions of the effective Hamiltonian in Eq. (B6) are invariant under a Fourier transform, up to a constant prefactor $(-i)^m$.

APPENDIX C: ENVELOPE FUNCTION APPROXIMATION—DERIVATION OF THE APPROXIMATE GRID STATES

In the following, we derive the approximate grid states introduced in Sec. IV B, by using the envelope function approximation. For convenience, we rescale the Hamiltonian in Eq. (30) by $\hbar\omega_0^2/\omega_c$, leading to

$$H = \frac{P^2 + X^2}{2} - W[\cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)], \quad (\text{C1})$$

where we introduce the large dimensionless parameter $W = \omega_c V_0 / \hbar\omega_0^2 \gg 1$.

Let us first examine the symmetries of this Hamiltonian.

First, H is an even function of both X and P , and so, in both X and P representation, its eigenfunctions can be chosen to be even and odd real-valued functions.

Second, since the Hamiltonian is invariant under an exchange of the position and momentum variables ($X \leftrightarrow P$), its eigenfunctions must be equal (up to an overall phase) in both representations. From this statement, it follows that the nondegenerate eigenfunctions of the Hamiltonian must be eigenfunctions of the Fourier transform.

In general, applying the Fourier transform twice is equivalent to applying a parity operation; applying the Fourier transform four times corresponds to the identity. For this reason, the eigenvalues of the Fourier transform are integer powers of i . In particular, the eigenfunctions of the Fourier transform with even parity have eigenvalues ± 1 under Fourier transform.

After having analyzed the Hamiltonian's symmetries, we now determine its approximate eigenfunctions with a consequent double application of the envelope function approximation [66,82,83].

Because $W \gg 1$, we start from the GKP Hamiltonian

$$H_{\text{GKP}} = -W[\cos(2\sqrt{\pi}X) + \cos(2\sqrt{\pi}P)], \quad (\text{C2})$$

and we treat the extra terms as smooth perturbations, determining the behavior of the envelope function that modulates the GKP ground states. The eigenstates of H_{GKP} are uniquely defined by the Zak states [28–30,126]

$$e^{i2\sqrt{\pi}X}|\Psi_{k,q}\rangle = e^{i2\sqrt{\pi}q}|\Psi_{k,q}\rangle, \quad (\text{C3a})$$

$$e^{i\sqrt{\pi}P}|\Psi_{k,q}\rangle = e^{i\sqrt{\pi}k}|\Psi_{k,q}\rangle, \quad (\text{C3b})$$

with $k \in [-\sqrt{\pi}/2, 3\sqrt{\pi}/2)$ and $q \in [-\sqrt{\pi}/2, \sqrt{\pi}/2)$. In position representation, these states can be written in the Bloch form

$$\Psi_{k,q}(X) = e^{ikX} \sum_{n \in \mathbb{Z}} \delta(X - \sqrt{\pi}n - q), \quad (\text{C4})$$

and, by a Fourier transform, we obtain (up to a global prefactor)

$$\Psi_{k,q}(P) = e^{-iqP} \sum_{n \in \mathbb{Z}} \delta(P - 2\sqrt{\pi}n - k). \quad (\text{C5})$$

The spectrum of H_{GKP} is continuous, and its band structure is given by

$$E(k, q) = -W[\cos(2\sqrt{\pi}k) + \cos(2\sqrt{\pi}q)]. \quad (\text{C6})$$

This band has two degenerate minima with energy $E = -2W$, obtained for $(k = 0, q = 0)$ and $(k = \sqrt{\pi}, q = 0)$.

Let us now consider the Hamiltonian $H_P = H_{\text{GKP}} + P^2/2$. Because the latter term is smooth on the scale of the GKP Hamiltonian, we assume that the eigenfunction of H_P can be factorized as

$$\tilde{\Psi}_{k,q}(P) = \phi(P)\Psi_{k,q}(P), \quad (\text{C7})$$

where the Bloch function $\Psi_{k,q}(P)$ is given in Eq. (C5). The function $\phi(P)$ is a smooth envelope that modulates the Bloch function, and, in analogy to solid-state theory, it satisfies [66,82,83]

$$\left(E(k, q) + \frac{P^2}{2}\right)\phi(P) = \mathcal{E}(k)\phi(P), \quad (\text{C8})$$

with the eigenvalue $\mathcal{E}(k)$. We are interested in the ground state eigenfunctions only, and so we expand $E(k, q)$ around the minimum $q = 0$ (effective mass approximation). Neglecting a constant energy offset and promoting the crystal momentum q to the operator $i\partial_P$, we obtain the harmonic oscillator differential equation

$$\left(-2\pi W\partial_P^2 + \frac{P^2}{2}\right)\phi(P) = [\mathcal{E}(k) + W\cos(2\sqrt{\pi}k)]\phi(P), \quad (\text{C9})$$

which has the ground state wave function

$$\phi(P) = \frac{\sqrt{\Delta}}{\pi^{1/4}} \exp\left(-\frac{\Delta^2 P^2}{2}\right). \quad (\text{C10})$$

The characteristic length of this oscillator is

$$1/\Delta = (4\pi W)^{1/4} \quad (\text{C11})$$

and corresponds to the broadening $1/\Delta$ of the Gaussian envelope function discussed in Sec. IV B; see Eq. (33).

To include the $X^2/2$ term, we first Fourier transform Eq. (C7) for $q = 0$, leading to the Bloch functions in X :

$$\tilde{\Psi}_{k,0}(X) \approx e^{ikX} \sum_{n \in \mathbb{Z}} e^{-(X - \sqrt{\pi}n + k)^2/2\Delta^2}, \quad (\text{C12})$$

where the approximate sign holds in the limit $\Delta \ll 1$. Because $X^2/2$ varies smoothly in each period of the Bloch function, we proceed as before and factorize the wave function of $H = H_P + X^2/2$ as

$$\psi(X) = \Phi(X)\tilde{\Psi}_{k,0}(X), \quad (\text{C13})$$

where the modulating function $\Phi(X)$ satisfies the eigenvalue equation

$$\left(\mathcal{E}(k) + \frac{X^2}{2}\right)\Phi(X) = \epsilon\Phi(X) \quad (\text{C14})$$

and, in the effective mass approximation, is given by

$$\Phi(X) = \frac{\sqrt{\Delta}}{\pi^{1/4}} \exp\left(-\frac{\Delta^2 X^2}{2}\right). \quad (\text{C15})$$

Importantly, $\mathcal{E}(k)$ has two degenerate minima at $k = 0$ and at $k = \sqrt{\pi}$, and so we obtain two approximate ground state eigenfunctions, that are given by the broadened GKP code words $\psi_0(X)$ and $\psi_1(X)$ defined in Eq. (32). Note that these states are approximately orthonormal in the limit $\Delta \ll 1$. Importantly, in the approximation used here, the eigenstates are degenerate, and the Fourier transforms of $\psi_{0,1}(X)$ are approximately given by

$$\psi_{\pm}(X) = \frac{\psi_0(X) \pm \psi_1(X)}{\sqrt{2}}. \quad (\text{C16})$$

Thus, we construct the states $\psi_{H\pm}(X)$ [see Eq. (31)] that are approximately even and odd functions under Fourier transform and respect the symmetries of the Hamiltonian.

As a consistency check, we can also estimate the broadening of the code words in Fig. 7, by using the Heisenberg uncertainty principle. For weak confinements $\hbar\omega_0^2/\omega_c V_0 \ll 1$, we expect the low-energy eigenfunctions of the Hamiltonian in Eq. (30) to have support only in the vicinity of $X = n2\sqrt{\pi}$ and $X = n2\sqrt{\pi} + \sqrt{\pi}$, respectively, with $n \in \mathbb{Z}$.

Let us focus on what happens close to $X = 0$. From Heisenberg's uncertainty principle, confining a particle to a narrow region causes large fluctuations in momentum P , such that $\hbar\omega_0^2\langle P^2 \rangle/\omega_c \gg V_0$. By expanding $\cos(2\sqrt{\pi}X)$ up to second order and neglecting the fast-oscillating term $\cos(2\sqrt{\pi}P)$ and the small perturbation $\hbar\omega_0^2 X^2/\omega_c$, we obtain

$$H \approx \frac{\hbar\omega_0^2 P^2}{2\omega_c} + 2\pi V_0 X^2. \quad (\text{C17})$$

This Hamiltonian is the one of a harmonic oscillator, and its ground state is a Gaussian of width

$$\Delta = \sqrt[4]{\frac{\hbar\omega_0^2}{4\pi\omega_c V_0}}. \quad (\text{C18})$$

This wave function approximates the narrow Gaussian centered at $X = 0$ of the approximate GKP state $\psi_0(X)$.

The width of the wide Gaussian envelope can be found by a similar argument, where one first considers localization in momentum P and then Fourier transforms the result, leading to Eq. (C11).

APPENDIX D: DERIVATION OF THE INTEGRATION KERNEL

In this Appendix, we derive the analytical expression of the integration kernel in Eq. (37), connecting the wave functions of the two-dimensional Hamiltonians and that of the one-dimensional Hamiltonians projected onto the LLL. By inserting the identity operator

$$\hat{\mathbf{1}} = \sum_{n=0}^{\infty} \int_{-\infty}^{\infty} dX |n, X\rangle \langle n, X|, \quad (\text{D1})$$

in $\Psi(x_1, x_2) = \langle x_1, x_2 | \Psi \rangle$, and neglecting mixing to higher Landau levels, we find

$$K_0(x_1, x_2; X) = \langle x_1, x_2 | 0, X \rangle. \quad (\text{D2})$$

To derive Eq. (37), we introduce the annihilation operators of the dynamical momenta and guiding center variables:

$$\hat{a} = \frac{1}{\sqrt{2}} \frac{l_B}{\hbar} (\hat{\pi}_2 + i\hat{\pi}_1), \quad \hat{b} = \frac{1}{\sqrt{2}} \frac{1}{l_B} (\hat{R}_1 + i\hat{R}_2). \quad (\text{D3})$$

Note that, for the sake of clarity, within this Appendix, we indicate operators with hats on top. Also, from now on, the coordinates x_i are given in magnetic units, thus scaled by the magnetic length l_B . Inserting the definition of the guiding center variables $\hat{R}_i$ [cf. Eq. (12)] in this equation yields

$$\hat{b} = \frac{\hat{x}_1 + i\hat{x}_2}{\sqrt{2}} - \hat{a}^\dagger. \quad (\text{D4})$$

The projector $\hat{\Pi}_{\text{LLL}} = |0\rangle\langle 0|_a$ onto the LLL acts solely in the Hilbert space of the dynamical momenta and, thus, commutes with any operator acting exclusively on the guiding center variables. Because $[\hat{\Pi}_{\text{LLL}}, \hat{b}] = 0$, we find that the action of $\hat{b}$ on the projected position eigenstate,

$$\hat{b}\hat{\Pi}_{\text{LLL}}|x_1, x_2\rangle = \frac{x_1 + ix_2}{\sqrt{2}} \hat{\Pi}_{\text{LLL}}|x_1, x_2\rangle, \quad (\text{D5})$$

is equivalent to the definition of a coherent state [56] $|\beta\rangle_b$ in the subspace of the guiding center variables, i.e.,

$$\hat{\Pi}_{\text{LLL}}|x_1, x_2\rangle \propto |0\rangle_a \otimes |\beta = (x_1 + ix_2)/\sqrt{2}\rangle_b. \quad (\text{D6})$$

We stress that Eq. (D5) is valid only for projectors onto the *lowest* Landau level, since $\hat{\Pi}_{\text{LLL}}\hat{a}^\dagger = 0$.

In conclusion, the integration kernel as defined in Eq. (D2) is the complex conjugate of the coherent state wave function in position representation [56]:

$$K_0(x_1, x_2; X) = \frac{1}{\sqrt{2\pi^{3/4}}} \exp\left(-\frac{(X - x_1)^2}{2}\right) \times \exp(-ix_2 X) \exp[ig(x_1, x_2)], \quad (\text{D7})$$

where the prefactor is chosen to satisfy

$$\iint dx_1 dx_2 K_0^*(x_1, x_2; X) K_0(x_1, x_2; X') = \delta(X - X'). \quad (\text{D8})$$

The real-valued function $g(x_1, x_2)$ arranges an adjustment of the global complex phase for fixed values of x_i . It is determined by the chosen gauge of the vector potential $A(x_1, x_2)$ and is fixed by demanding the state $|0, X\rangle$ to lie in the LLL. Since the annihilation operator $\hat{a}$ of the dynamical momenta, expressed in the initial positions $\hat{x}_i$, is gauge dependent, the integration kernel has to satisfy (in magnetic units)

$$0 = [\partial_{x_1} - i\partial_{x_2} + A_2(x_1, x_2) + iA_1(x_1, x_2)]K_0(x_1, x_2; X), \quad (\text{D9})$$

leading to the equation

$$-\nabla g(x_1, x_2) = \begin{pmatrix} A_1(x_1, x_2) \\ A_2(x_1, x_2) - x_1 \end{pmatrix} \quad (\text{D10})$$

for the function $g(x_1, x_2)$. In particular, for the symmetric gauge (in magnetic units),

$$A_1(x_1, x_2) = -\frac{x_2}{2}, \quad A_2(x_1, x_2) = \frac{x_1}{2}, \quad (\text{D11})$$

we find, up to a trivial constant,

$$g(x_1, x_2) = \frac{x_1 x_2}{2}. \quad (\text{D12})$$

Combining Eqs. (D7) and (D12), we obtain Eq. (37).

APPENDIX E: RELATION BETWEEN THE FOURFOLD ROTATION AND THE FOURIER TRANSFORM

Given the integral transform in Eq. (36), here, we want to show that any one-dimensional eigenfunction of the Fourier transform with eigenvalue $(-i)^n$, i.e.,

$$\frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} dP \psi(P) e^{-iXP} = (-i)^n \psi(X), \quad (\text{E1})$$

results in a two-dimensional eigenfunction of the fourfold rotation with the complex conjugate eigenvalue. To this end, we make use of the result for the integration kernel in symmetric gauge [see Eq. (37)] and obtain

$$\begin{aligned} \Psi(x_2, -x_1) &= \int_{-\infty}^{\infty} dX K_0(x_2, -x_1; X) \psi(X) \\ &= \frac{i^n}{2\pi^{7/4}} \int_{\mathbb{R}^2} dX dP e^{-(X-x_2)^2/2} e^{i(x_1-P)X} e^{-x_1 x_2/2} \psi(P) \\ &= \frac{i^n}{\sqrt{2\pi^{3/4}}} \int_{-\infty}^{\infty} dP e^{-(P-x_1)^2/2} e^{-ix_2 P} e^{ix_1 x_2/2} \psi(P) \\ &= i^n \int_{-\infty}^{\infty} dP K_0(x_1, x_2; P) \psi(P) \\ &= i^n \Psi(x_1, x_2). \end{aligned} \quad (\text{E2})$$

The particular choice of the symmetric gauge is essential for the previous derivation, since it determines the complex phase of the integration kernel for fixed values of x_i . This choice is in agreement with the observation that the two-dimensional Hamiltonian in Eq. (28) is fourfold rotational symmetric in the $x_1 x_2$ plane in the symmetric gauge only. Rotation-symmetric bosonic codes are analyzed in Ref. [127].

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