Confinement and Lattice Quantum-Electrodynamic Electric Flux Tubes Simulated with Ultracold Atoms

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> We propose a method for simulating $(2 + 1)D$ compact lattice quantum-electrodynamics, using ultracold atoms in optical lattices. In our model local Bose-Einstein condensates' (BECs) phases correspond to the electromagnetic vector potential, and the local number operators represent the conjugate electric field. The well-known gauge-invariant Kogut-Susskind Hamiltonian is obtained as an effective low-energy theory. The field is then coupled to external static charges. We show that in the strong coupling limit this gives rise to ''electric flux tubes'' and to confinement. This can be observed by measuring the local density deviations of the BECs, and is expected to hold even, to some extent, outside the perturbative calculable regime.

> DOI: [10.1103/PhysRevLett.107.275301](http://dx.doi.org/10.1103/PhysRevLett.107.275301) PACS numbers: 67.85.Hj, 11.15.Ha

Free quarks are not found in nature. This is due to the mechanism of confinement. A lot of theoretical progress in this area has been achieved—either in the lattice Euclidean approach by Wilson [[1\]](#page-3-0), in nonperturbative methods by Polyakov [\[2](#page-3-1)], or using the lattice Hamiltonian formalism, by Kogut and Susskind [[3](#page-3-2)[,4\]](#page-3-3).

Although gauge theories can be latticized either in a compact (nonlinear) or noncompact (linear) manner, the compactness is essential to the confinement mechanism [\[5\]](#page-3-4). It has been shown that in an Abelian $3 + 1$ compact lattice gauge theory, a phase transition is supposed to take place between two phases—the Coulomb phase for small couplings, which exhibits the "regular" $V(R) \propto 1/R$ static potential between two R-separated static charges, and the confining phase, for which the static potential is linear in the distance between the charges— $V(R) \propto R$, for large values of the coupling constant [\[4](#page-3-3)]. (Non-Abelian theories, on the other hand, confine for all values of the coupling constant.) However, for an Abelian $2 + 1$ compact lattice gauge theory, confinement was shown to take place for all the values of the coupling constant, due to nonperturbative effects of instantons [\[2,](#page-3-1)[5](#page-3-4)–[7\]](#page-3-5). Considering thermal effects as well, even in $2 + 1$ dimensions a phase transition to a Coulomb phase exists for $T > 0$ [[8](#page-3-6)[,9\]](#page-3-7).

The mechanism responsible for confinement is believed to produce an ''electric flux tube,'' connecting two static charges in the confining phase, which is hard to measure directly. It requires measuring the force and/or potential between two static charges. If one wishes to observe the phase transition, the coupling constant has to be varied, which poses another difficulty. A quantum simulation of such a model could allow a direct test of the confinement mechanism and the phase transitions.

Quantum gases of ultracold atoms, implemented in optical lattices [[10](#page-3-8)], provide models with highly controllable parameters and offer a natural playground for the simulations of such models. Quantum simulation approaches of various kinds and aspects of compact $U(1)$ pure gauge theory, in cold gases and other systems, have been proposed by several authors: In [\[11\]](#page-3-9), an effective theory of $U(1)$ spin liquid in pyrochlore was discussed; in [\[12\]](#page-3-10), using a molecular state in optical lattices, an effective theory of ring exchange was derived, and it is, in the limit of no hopping, a U(1) lattice gauge theory, with a Coulomb phase; in [[13\]](#page-3-11), emergence of ''artificial photons'' and a Coulomb phase in an effective theory based on dipolar bosons in an optical lattice were shown; and in [\[14\]](#page-3-12), a possibility to simulate a spin U(1) pure gauge theory as a low-energy theory with a system of Rydberg atoms was presented.

In this Letter, we suggest a method for simulating compact QED with cold atoms in optical lattices, which should enable a direct observation of electric flux tubes that emerge in the mechanism of confinement. In our model, the vector potential and its conjugate electric field are represented by the local condensate phase operators and their conjugate number operators. These observables ''live'' on the links of a two- or three-dimensional optical lattice, and hence each link of the lattice is here represented by a separate Bose-Einstein condensate (BEC). In order to obtain the QED Hamiltonian, one has to generate certain two- and four-body interactions between the condensates, that manifest local gauge invariance. In order to avoid the hopping processes of an ordinary Bose-Hubbard model, we introduce a four-species two-dimensional setup (Fig. [1\)](#page-1-0). The condensates are located on the links of a lattice—each species on a different link—and overlap at the lattice's vertices. Hence, condensates of the same type are spatially separated, as depicted in Fig. [1,](#page-1-0) causing the attenuation of hopping processes. Next we use Raman transitions and two atom scattering processes in order to create special "diagonal" hopping and nonlinear interactions. As we show, in this new setup a certain choice of parameters gives rise to gauge invariance in the low-energy sector; hence compact QED emerges as an effective theory.

FIG. 1 (color online). Left: Structure of the lattice. The different condensate species are colored in four colors; the colored boxes represent the links (condensates), and there the localized wave functions are concentrated. At the vertex (symbolized by a cube) the wave functions of the neighboring links overlap and these are the only overlap integrals which are not negligible. Right: A close-up picture of a single vertex, showing the various interaction parts of the Hamiltonian—scattering and hopping.

To study the effect of confinement within this setup, we can introduce two spatially separated effective ''charges'' by creating local deformations of the trapping potential at the position of the charges at the relevant vertices. We then expect that the local atomic densities, within the QED parameter regime, should manifest the effect of confinement by the appearance of a fluxlike tube of alternating atomic density deviations along the line connecting the charges (Fig. [2\)](#page-1-1), while such a flux tube will not appear outside the QED parameters regime. Other possible implications of our model will be briefly discussed in the summary.

We begin with a system of condensates described by the Hamiltonian $H = \int d^3x \sum_{i,j=1}^4 \mathcal{H}_{ij}(\mathbf{x})$, where

$$
\mathcal{H}_{ij}(\mathbf{x}) = \Psi_i^{\dagger}(\mathbf{x})(\delta_{ij}(\mathcal{H}_0^i(\mathbf{x}) + V_M(\mathbf{x})) + \Omega_{ij})\Psi_j(\mathbf{x}) \n+ \frac{g_{ij}}{2}\Psi_i^{\dagger}(\mathbf{x})\Psi_j^{\dagger}(\mathbf{x})\Psi_j(\mathbf{x})\Psi_i(\mathbf{x}).
$$
\n(1)

 δ_{ij} is Kronecker's delta, g_{ij} are the s-wave scattering coefficients, and Ω_{ij} are Rabi frequencies. It contains the following parts. The "free" Hamiltonian of each species:

FIG. 2 (color online). An example of the charge and flux configurations, for $R = 2$. The different colors represent the condensate species. The upper couple of charges are with QED quantum numbers, and the lower couple with BEC local number deviations quantum numbers. Such a flux tube can be embedded, in the absence of other charges, in a lattice whose other links carry $E = 0$.

 $\mathcal{H}_0^i(\mathbf{x}) = -\frac{\nabla^2}{2m} + V_i(\mathbf{x})$, where $V_i(\mathbf{x})$ is the optical lattice
transition potential of the species *i*; the scattering terms, set trapping potential of the species i ; the scattering terms, set by the coupling constants g_{ij} (neglecting the three- and four-body interactions): (i) self-scattering terms, $g_{ii} \equiv g_1$,
(ii) two-species scattering terms $g_{i2} \equiv g_{2i} \equiv g_{2i} \equiv$ (ii) two-species scattering terms, $g_{12} = g_{21} = g_{34}$ $g_{43} \equiv g_2$ along straight lines, and along the diagonals,
 $g_{12} \equiv g_{21} \equiv g_{11} \equiv g_{22} \equiv g_{22} \equiv g_{22} \equiv g_{23} \equiv g_{33} \$ $g_{13} = g_{31} = g_{14} = g_{41} = g_{23} = g_{32} = g_{24} = g_{42} \equiv g_3$
(all the other g₁'s are zero); an "external charges" simu-(all the other g_{ii} 's are zero); an "external charges" simulating potential, which deforms the lattice potential at the vertices and is approximated by a very localized potential, $V_M(\mathbf{x}) \equiv \sum_{m,n} \alpha_{m,n} \delta(\mathbf{x} - \mathbf{x}_{m,n})$, where $\alpha_{m,n}$ are constants
(whose value and sign are related to the external charges (whose value and sign are related to the external charges and will be determined in the sequel) and $\mathbf{x}_{m,n}$ is the position of the (m, n) vertex. The laser generated Rabi terms $\Omega_{13} = \Omega_{31} = \Omega_{14} = \Omega_{41} = \Omega_{23} = \Omega_{32} = \Omega_{24} = \Omega_{41} = \Omega_{41} = \Omega_{42} = \Omega_{43} = \Omega_{44} = \Omega_{45} = \Omega_{46} = \Omega_{47} = \Omega_{48} = \Omega_{49} = \Omega_{40} = \Omega_{4$ $\Omega_{42} \equiv \Omega_0^7$ couple the condensates to each other in a special diagonal manner as depicted in Fig. 1. (All the other cial, diagonal manner, as depicted in Fig. [1](#page-1-0). (All the other Ω_{ij} 's are zero.) Since the minima of the same species are far enough apart, the hopping effects are solely controlled by the latter Rabi terms. Experimentally, our scheme can be implemented by using holographic masks techniques [\[15\]](#page-3-13) in order to generate the required optical lattice and using optical Feshbach resonances in order to control the coupling strengths g_{ij} [\[16](#page-3-14)[–18\]](#page-3-15). Raman transitions can be used to control the coefficients Ω_{ij} of the Rabi terms.

The second quantization wave functions of the condensates (taking into account only the lowest band excitations) are $\Psi_{1,2} = \sum_{m,n} a_{m,n} \phi_{m,n}(\mathbf{x}), \quad \Psi_{3,4} = \sum_{m,n} b_{m,n} \chi_{m,n}(\mathbf{x}),$
where a b are single-mode annihilation operators where $a_{m,n}$, $b_{m,n}$ are single-mode annihilation operators, annihilating one particle in the ground state of the corresponding link (minimum). Note that because of the lattice's structure, not all the values (m, n) are included in the wave function of each species. We assume that the local Wannier functions [[10](#page-3-8)] respect the symmetries $\phi_{m,n}(\mathbf{x}) =$ $\phi(\mathbf{x} - \mathbf{x}_{m,n}^{1,2}), \chi_{m,n}(\mathbf{x}) = \chi(\mathbf{x} - \mathbf{x}_{m,n}^{3,4}) = \phi(R\mathbf{x})$, where R is the appropriate rotation operator, and that they are chosen to be real [\[19\]](#page-3-16).

Plugging the wave functions into the Hamiltonian ([1\)](#page-1-2), one gets, using the above assumptions, that the only nonnegligible contributions are $2\lambda + \mu = \frac{g_1}{2} \int d^3x |\dot{\phi}(\mathbf{x} - \mathbf{y})|^4 = \frac{g_2}{2} \int d^3x |\dot{\phi}(\mathbf{x} - \mathbf{y})|^4$ $\langle \mathbf{x}_0 |^{4}, \quad V_0 \equiv \int d^3x \phi^*(\mathbf{x} - \mathbf{x}_0) \left[-\frac{\nabla^2}{2m} + V_i(\mathbf{x}) \right] \phi(\mathbf{x} - \mathbf{x}_0),$
 $V = \int d^3x \phi(\mathbf{x} - \mathbf{x}_0) \left[\frac{2}{3m} \phi(\mathbf{x} - \mathbf{x}_0) \right] \phi(\mathbf{x} - \mathbf{x}_0).$ $V_2 = \int d^3x |\phi(\mathbf{x} - \mathbf{x}_0)|^2 |\phi(\mathbf{x} - \mathbf{x}_1)|^2$, $V_3 = \int d^3x |\phi(\mathbf{x} - \mathbf{x}_0)|^2 |\phi(\mathbf{x} - \mathbf{x}_0)|^2$ $|\mathbf{x}_0|^2 |\chi(\mathbf{x} - \mathbf{x}_2)|^2$, $\Delta_{m,n} \equiv -\frac{1}{2\lambda} \alpha_{m,n} |\phi(\mathbf{x}_{m,n})|^2$, and $\Omega_0 \equiv \Omega_0^{\prime} \int d^3x \phi^*(\mathbf{x} - \mathbf{x}_0) \chi(\mathbf{x} - \mathbf{x}_2)$ (here the reality of the Wannier functions is employed). \mathbf{x}_0 is the position of an Wannier functions is employed). x_0 is the position of an arbitrary minimum of the potential (due to the symmetries), x_1 is an adjacent minimum in the same direction (separated by a single lattice spacing), and x_2 is an adjacent minimum in the orthogonal direction (rotated). In the following we assume that g_2, g_3 satisfy the relation $g_2V_2 = g_3V_3 = 2\lambda.$

Let $N_{m,n}^k$ be the local number operators, emanating from the vertex (m, n) : for horizontal (\hat{x}) links $k = 1$ and for vertical ($\hat{\mathbf{y}}$) links $k = 2$. $N_T = \sum_{m,n,k} N_{m,n}^k$, the total number

of particles, is a constant of motion. We choose a subspace by fixing $N_T = \mathcal{N}_L N_0$, where \mathcal{N}_L is the number of links and $N_0 \gg 1$. Defining $M_{m,n} = 4N_0 + \Delta_{m,n}$, $G_{m,n} =$ $N_{m,n}^1 + N_{m,n}^2 + N_{m-1,n}^1 + N_{m,n-1}^2 - M_{m,n}$, after some algebra one obtains the Hamiltonian $H = H - H$ bra, one obtains the Hamiltonian $H_0 = H - H_R =$
 $\lambda \sum_{i=1}^{N} G_i^2 + \mu \sum_{j=1}^{N} (N^k)^2$. The nearest-peighbor hop- $\lambda \sum_{m,n} G_{m,n}^2 + \mu \sum_{m,n,k} (N_{m,n}^k)^2$. The nearest-neighbor hop-
ping part, which results here from the Rabi terms, can be $\lambda_{\sum_{m,n}} G_{m,n}^T + \mu_{\sum_{m,n,k}} (N_{m,n})^T$. The heartst-heighbor hop-
ping part, which results here from the Rabi terms, can be written as $H_R = \Omega_0 \sum_{m,n} (a_{m,n} b_{m,n}^\dagger + a_{m,n} b_{m+1,n}^\dagger +$ $a_{m,n+1}b_{m,n}^{\dagger} + a_{m,n+1}b_{m+1,n}^{\dagger} + \text{H.c.}.$

Gauss's law.—We wish to obtain a gauge-invariant theory, and hence we would like to constrain Gauss's law on the system. This is satisfied in the QED regime: $\lambda \gg \mu$ and $\lambda \gg \Omega_0$, in which H_R can be treated as a small
perturbation. Let us first find the ground state of H_0 . perturbation. Let us first find the ground state of H_0 . After expanding the number operator on each link around $N_0, N_{m,n}^k = N_0 + \delta_{m,n}^k$, one obtains at each vertex $G_{m,n} =$
 $S^1 + S^2 + S^1 + S^2 = \Lambda$ Within the sub- $\delta_{m,n}^1 + \delta_{m,n}^2 + \delta_{m-1,n}^1 + \delta_{m,n-1}^2 - \Delta_{m,n}$. Within the subspace of a constant, conserved N_T , $\sum_{m,n,k} \delta_{m,n}^k = 0$.
Neglecting constants of motion, one can rewrite the Neglecting constants of motion, one can rewrite the Hamiltonian in terms of $\delta_{m,n}^k$: $H_0 = \lambda \sum_{m,n} G_n^2$
 $\mu \sum (\delta^k)_{n=1}^k H_0 + H_0 \lambda \gg \mu$ and hence manntonian in terms or $\sigma_{m,n}^*$. $H_0 - \lambda \sum_{m,n} \sigma_{m,n}^* \mp \mu \sum_{m,n,k} (\delta_{m,n}^k)^2 \equiv H_G + H_E$. $\lambda \gg \mu$, and hence one
would like to minimize H₋ first. Thus we get that in the $\mu_{\sum_{m,n,k}}(o_{m,n})^2 = n_G + n_E$. $\lambda \gg \mu$, and hence one would like to minimize H_G first. Thus we get that in the ground state the sum of $\delta_{m,n}^k$'s around each vertex equals the $\Delta_{m,n}$ of the vertex: This imposes a modified Gauss's law (sum instead of discrete divergence), and hence the $\Delta_{m,n}$'s must be integers (positive, zero, or negative)—this can be set by adjusting the values of the $\alpha_{m,n}$'s in $V_M(\mathbf{x})$.
Next, to minimize the entire H_c (including H_c) we would Next, to minimize the entire H_0 (including H_E) we would like to choose the lowest $\delta_{m,n}^k$'s which satisfy this constraint.

Define the sublattices $A = \{(m, n) : m + n = \text{even}\},\$ $B = \{(m, n): m + n = \text{odd}\}\.$ Note that for states that respect Gauss's law (for which $G_{m,n}|\psi\rangle = 0$), which will later be the physically interesting states, the sum of $\Delta_{m,n}$'s of each sublattice must be zero. This follows from adding the $G_{m,n}$'s of each sublattice, taking into account that the total particle number deviation is zero.

Quantum rotor approximation.—If we set that at each vertex $|\Delta_{m,n}| \ll N_0$, we get that on each link, in the ground state of H_0 , within our subspace, $|\langle \delta_{m,n}^k \rangle| \ll N_0$, and thus, after taking into account the perturbative corrections, one obtains that on each link $\sigma_{m,n}^k \equiv \langle (\delta_{m,n}^k)^2 \rangle^{1/2} \ll N_0$. Note botants that on each link $\sigma_{m,n} = \langle (\sigma_{m,n})^2 \rangle^3 \ll N_0$. Note
that $\sigma_{m,n}^k \ll \frac{\lambda N_0 (N_0 - 1)}{\Omega_0}$ (because $\frac{\Omega_0}{\lambda} \ll N_0$), and hence the two conditions of [[20](#page-3-17)] are fulfilled and the Hamiltonian can be approximated as a quantum rotor Hamiltonian. H_G , H_E remain the same because they are already written in the number deviations' notation. Because of the phase-number relation of the condensates, $[N_{m,n}^i, \theta_{m',n'}^j] = i \delta_{mm'} \delta_{ij},$ $[N_{m,n}^i, e^{\pm i\theta_{m,n}^i}] = \pm e^{\pm i\theta_{m,n}^i}$, and therefore we can define phase-only lowering and raising operators, $\tilde{a}_{m,n} = e^{i\theta_{m,n}^1}$
 $\tilde{a}^{\dagger} = e^{-i\theta_{m,n}^1}$, $\tilde{b} = e^{i\theta_{m,n}^2}$, $\tilde{b}^{\dagger} = e^{-i\theta_{m,n}^2}$, which operate $\tilde{a}^{\dagger}_{m,n} = e^{-i\theta^1_{m,n}}, \tilde{b}_{m,n} = e^{i\theta^2_{m,n}}, \tilde{b}^{\dagger}_{m,n} = e^{-i\theta^2_{m,n}},$ which operate
on the local number deviations: \tilde{a} $|\tilde{\lambda}^1|$ $\rangle = |\tilde{\lambda}^1| - 1$ on the local number deviations: $\tilde{a}_{m,n} |\delta_{m,n}^1\rangle = |\delta_{m,n}^1 - 1\rangle$, etc., and since $N_0 \gg 1$, $\sqrt{N_0(N_0 + 1)} \approx N_0$. Thus one gets $H_R = \Omega \sum_{m,n} (\tilde{a}_{m,n} \tilde{b}_{m,n}^{\dagger} + \tilde{a}_{m,n} \tilde{b}_{m+1,n}^{\dagger} + \tilde{a}_{m,n+1} \tilde{b}_{m,n}^{\dagger} +$
 $\tilde{a}_{m,n+1} \tilde{b}_{m,n+1}^{\dagger}$ $\tilde{a}_{m,n+1}\tilde{b}_{m+1,n}^{\dagger} + \text{H.c.}$, where $\Omega = \Omega_0 N_0$.

Effective Hamiltonian. Let us look a

Effective Hamiltonian.—Let us look again at the eigenstates and eigenvalues of H_G . Since $[H_G, H_E] = 0$, the two Hamiltonians can be mutually diagonlized. The eigenstates of H_E are number states, and we shall use this basis to diagonalize H_G as well. Since $\lambda \gg \mu$, Ω , the Gauss
Hamiltonian H_G is much stronger than the other two and Hamiltonian H_G is much stronger than the other two, and therefore one can obtain an effective low-energy theory perturbatively [[21](#page-3-18)]. It is physically reasonable to derive an effective Hamiltonian by projecting to the ground state manifold of H_G . Let us denote this manifold by $M: M =$ $\{ |M_{\alpha}\rangle : H_G|M_{\alpha}\rangle = 0 \}$. One can see that it is the physical subspace of states which respects Gauss's law. The pertursubspace of states which respects Gauss's law. The perturbative expansion to second order leads to $H_{\text{eff}} = H_E +$ H_B , where $H_B = -\frac{2\Omega^2}{\lambda}$ $\frac{\Omega^2}{\lambda} \sum_{m,n} (\tilde{a}^{\dagger}_{m,n+1} \tilde{b}_{m,n} \tilde{a}^{\dagger}_{m,n} \tilde{b}_{m+1,n} + \text{H.c.})$ is the desired gauge-invariant four-body plaquette interaction.

Compact QED analogy.—We next relate this model to compact QED and discuss the implications. First, let us switch to QED-like variables. In order to do so, we perform the transformation $E_{m,n}^k \equiv (-1)^{m+n} \delta_{m,n}^k$, $Q_{m,n} \equiv (-1)^{m+n} \Delta$ and $\theta^k \to (-1)^{m+n} \theta^k$ Because of the $\bar{\theta}(-1)^{m+n}\Delta_{m,n}$, and $\theta_{m,n}^k \to (-1)^{m+n}\theta_{m,n}^k$. Because of the transformation of the phases of links emanating from subtransformation of the phases of links emanating from sublattice B vertices, these links' raising and lowering operators have to be swapped.

This transforms H_E , which can be identified as the "electric Hamiltonian," to $H_E = \mu \sum_{m,n,k} (E_{m,n}^k)^2$, electric rialimionian, to $H_E - \mu_{\sum_{m,n,k}}(E_{m,n})$
and $H_B = -\frac{2\Omega^2}{\lambda} \sum_{m,n} (\tilde{a}_{m,n}^\dagger \tilde{a}_{m,n+1} \tilde{b}_{m+1,n}^\dagger \tilde{b}_{m,n} + \text{H.c.})$ $\frac{\Omega^2}{\lambda} \sum_{m,n} (\tilde{a}_{m,n}^\dagger \tilde{a}_{m,n+1} \tilde{b}_{m+1,n}^\dagger \tilde{b}_{m,n} + \text{H.c.}) =$ $-\frac{4\Omega^2}{\lambda}$ $-\frac{4\Omega^2}{\lambda} \sum_{m,n} \cos(\theta_{m,n}^1 + \theta_{m+1,n}^2 - \theta_{m,n+1}^1 - \theta_{m,n}^2)$ is the magnetic part of the compact QED Hamiltonian (the cosine's argument is the discrete curl of $\theta_{m,n}$, which is the magnetic field). Thus we obtained an effective low-energy theory whose Hamiltonian is the compact QED Hamiltonian, constrained with Gauss's law (which is the low-energy constraint):

$$
G_{m,n}|\psi\rangle = (-1)^{m+n}(\text{div}\mathbf{E}_{m,n} - Q_{m,n})|\psi\rangle = 0. \qquad (2)
$$

Confinement of external static charges.—Define a new finite energy scale, $U_0 = \frac{2}{g^2} \mu = \frac{4\Omega^2 g^2}{\lambda}$, and rescale the Hamiltonian to

$$
\bar{H} = H_{\text{eff}}/U_0 = \frac{g^2}{2} \sum_{m,n,k} (E_{m,n}^k)^2
$$

$$
- \frac{1}{g^2} \sum_{m,n} \cos(\theta_{m,n}^1 + \theta_{m+1,n}^2 - \theta_{m,n+1}^1 - \theta_{m,n}^2), \qquad (3)
$$

which is the well-known Kogut-Susskind Hamiltonian for an Abelian lattice gauge theory [\[3,](#page-3-2)[4](#page-3-3)]. From the definition of U, one gets $g^4 = \frac{\lambda \mu}{2\Omega^2}$. This Hamiltonian has two limits: (i) The strong coupling limit, $g \gg 1$, or $\frac{\mu}{\Omega} \gg \frac{\Omega}{\lambda}$. In this limit, we can treat H_B as a perturbation to H_E . (ii) The weak coupling limit, $g \ll 1$, or $\frac{\mu}{\Omega} \ll \frac{\Omega}{\lambda}$. In this limit, we can treat H_E as a perturbation to H_B . In a 3 + 1 theory, the strong coupling limit is within the confining phase and the weak coupling limit is within the Coulomb phase, and a phase transition is expected in between $[1-5]$ $[1-5]$ $[1-5]$. In a $2 + 1$ theory, there is no phase transition and confinement is expected to occur for all $g > 0$ [\[5,](#page-3-4)[7\]](#page-3-5).

The external charges are limited by the restrictions imposed by the constraint $\sum_{(m,n)\in A}\Delta_{m,n}=0$, imposed by the constraint $\sum_{(m,n)\in A} \Delta_{m,n} = 0$,
 $\sum_{(m,n)\in B} \Delta_{m,n} = 0$. Subtracting the second constraint from the first, one gets $\sum_{m,n} Q_{m,n} = Q_{\text{tot}} = 0$. Thus, the total charge has to be zero. If we add these constraints we total charge has to be zero. If we add these constraints, we get another constraint, $\sum_{m,n} \Delta_{m,n} = 0$. This constraint does
not seem to have a OED analogy, but it has to be satisfied in not seem to have a QED analogy, but it has to be satisfied in our model.

Consider the case of a system with two unit external charges, in the strong coupling limit. Thus we seek for the ground state configurations of H_E and treat H_B as a perturbation. These charges must be of opposite signs, in order to satisfy the charge restrictions. For simplicity, we assume that the charges are fixed at the vertices (m, n) and $(m + R, n)$. If $R = 1$, the charges are fixed at two vertices of different sublattices, and hence have to have the same sign in terms of $\Delta_{m,n}$. This, however, does not satisfy the charge restrictions, and one has to add more charges to the system. This is true, in fact, for any odd R. Therefore, we shall consider only the case of an even R , for two charges in the system.

Denote $|R\rangle$ as the state of two such external charges,
= 1 Q = ϵ = -1 In the strong coupling limit it $Q_{m_0,n_0} = 1$, $Q_{m_0+R,n_0} = -1$. In the strong coupling limit, it can be written as a perturbative series, whose zeroth order term is $|R_0\rangle = \prod_{m_0 \le m \le m_0 + R} \tilde{a}^{\dagger}_{m,n} |\{0\}\rangle$. This corresponds to a flux tube from a positive charge to a negative one. Thus, in the strong limit, we get, indeed, the expected strong coupling linear behavior of the energy,

$$
\bar{E}(R) = \frac{1}{U_0} E(R) = \frac{g^2}{2} R + O(g^{-6}).
$$
 (4)

The effect can then be observed by measuring the local density deviations $\delta_{m,n}^k$, which are expected, in the leading order, to have a magnitude 1 and alternating signs between the two charges. An example for $R = 2$ can be seen in Fig. [2.](#page-1-1) When R is too large, the energy of H_G is smaller than the energy of such a flux tube, and then the low-energy theory breaks, and a flux tube is no longer the state of minimal energy. The low-energy picture holds as long as the flux tube length satisfies $R \ll \lambda/\mu$.

Outside the strong coupling regime, such perturbative calculations are no longer valid. However, in $2 + 1$ dimensions the confinement holds for all values of g . Hence, the effect should be seen experimentally even slightly outside the strong coupling limit, although not in the weak limit (in order to fit with the quantum rotor approximation).

It may also be possible to experimentally observe the effect of a finite temperature, $T > 0$, on the model, including a phase transition.

Extensions of the model.—In this Letter, we have shown a method to simulate compact QED using BECs in optical lattices, as a way to observe charge confinement. The suggested model can be extended in several ways. More realistic, nonperiodic boundary conditions can be imposed (e.g., no charges on the boundary). Using additional condensates (of new species), a $3 + 1$ simulation could be achieved. Interestingly, a dynamical charge which is minimally coupled to the field can be implemented using another condensate species. This is equivalent to a special case of the model of Fradkin and Shenker [[22](#page-3-19)], in which a Higgs field with a "frozen" radius is coupled to a $U(1)$ gauge field [\[23\]](#page-3-20).

To conclude, we hope that this model can serve as one of the building blocks of the bridge into the world of dynamic quantum gauge field theories simulations.

The authors would like to thank A. Casher, J. I. Cirac, S. Dürr, O. Kenneth, S. Nussinov, and B. Svetitsky for helpful discussions. This work has been supported by the Israel Science Foundation, the German-Israeli Foundation, and the European Commission (PICC).

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