Phase transition of light in circuit QED lattices coupled to nitrogen-vacancy centers in diamond

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We propose a hybrid quantum architecture for engineering a photonic Mott insulator-superfluid phase transition in a 2*D* square lattice of superconducting transmission line resonator (TLR) coupled to a single nitrogenvacancy (NV) center encircled by a persistent current qubit. The localization-delocalization transition results from the interplay between the on-site repulsion and the nonlocal tunneling. The phase boundary in the case of photon hopping with real-valued and complex-valued amplitudes can be obtained using the mean-field approach. Also, the quantum jump technique is employed to describe the phase diagram when the dissipative effects are considered. The unique feature of our architecture is the good tunability of effective on-site repulsion and photon-hopping rate, and the local statistical property of TLRs which can be analyzed readily using present microwave techniques. Our work opens new perspectives in quantum simulation of condensed-matter and many-body physics using a hybrid spin circuit QED system. The experimental challenges are realizable using current available technologies.

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I. INTRODUCTION

The microscopic properties of strongly correlated manyparticle systems emerging in solid-state physics are in general very hard to access experimentally^{[1,](#page-4-3)[2](#page-4-4)}. So how to simulate the properties of condensed-matter models using nontraditional controllable systems is highly desirable. Recently, the investigation of quantum simulation in the photon-based many-body physics has received much attention in different systems $¹⁻⁴$ $¹⁻⁴$ $¹⁻⁴$.</sup> Especially, there has been a great interest in mimicking quantum phase transition (QPT) of light with scalable coupled resonator array in the context of cavity/circuit quantum electrodynamics $(QED)^{5-8}$ $(QED)^{5-8}$ $(QED)^{5-8}$, which provides a convenient controllable platform for studying the strongly correlated states of light via photonic processes. On the other hand, the artificially engineered hybrid devices can permit measurement access with unique experimental control^{[10,](#page-4-8)[11](#page-4-9)}; and it is intriguing to employ a well-controllable quantum system with a tunable Hamiltonian to simulate the physics of another system of interest. This paradigm has promoted many experimental/theoretical proposals on probing the light phase and opened various possibilities for the simulation of many-body physics.

In this work, we develop an optical system for engineering the strongly correlated effects of light in a hybrid solid-state system. We consider a 2D square lattice of coupled $TLRs¹²$ $TLRs¹²$ $TLRs¹²$, where each TLR is coupled to a single $NV^{13,14}$ $NV^{13,14}$ $NV^{13,14}$ $NV^{13,14}$ encircled by a persistent current qubit (PCQ). We show that the competition between the NV-PCQ-TLR interaction and the nonlocal hopping induces the photonic localization-delocalization transition. Subsequently the Mott insulator (MI) phase and the superfluid (SF) phase can appear in a controllable way. The phase boundary in the case of photon hopping with real/complex-valued amplitudes can be obtained using the mean-field approach. Also, the quantum jump technique is employed to describe the phase diagram when the dissipation is considered. Finally, the possibility of observation of the

FIG. 1. (Color online) (a) Schematic circuit for the resonator lattice, where each TLR is coupled to a single NV center encircled by a PCQ, and the circles denote the central coupler. (b) The subsystem consisting of NV and PCQ, where NV is at the center of the loop. The PCQ is made up of three Josephson junctions, and it couples to the NV via the magnetic field at the center of the loop generated by the persistent currents in the loop. The energy diagram of the NV is shown in the red box.

QPT is discussed by employing experimentally accessible parameters.

In our architecture, one can tune independently the on-site emitter-field interaction and the nonlocal photonic hopping between adjacent TLRs. This permits us to systematically study the localization-delocalization transition of light in a complete parameter space. The main motivation for building such a hybrid system is to combine several advantages: *in situ* tunability of circuit elements^{[9](#page-4-13)}, spectroscopic technology for state readout, peculiar characteristics of NV (e.g., individual addressing and long coherence time at room-temperature^{[15](#page-4-14)}), and scalability of TLR arrays^{[6,](#page-4-15)[16](#page-4-16)[–18](#page-4-17)}. Recently, D. Underwood *et al* experimentally fabricated 25 arrays of TLRs and demonstrated the feasibility of quantum simulation in circuit QED system^{[19](#page-4-18)}. E. Lucero *et al* experimentally characterized a complex circuit composed of four phase qubits and five TLRs to realize intricate quantum algorithms^{[20](#page-4-19)}. The progress ren-

FIG. 2. (Color online) The dependence of effective on-site repulsion U_{+} on the photon number *n* under the different detunings Δ , where the solid (dotted) line denotes U_{+} (U_{-}).

ders the TLR lattice a good platform for studying condensedmatter physics with photons and makes our scheme to be more practical.

II. MODEL

As illustrated in Fig. 1, we consider a 2*D* lattice of coupled TLRs, where the basic unit consists of a TLR coupled to a single NV encircled by a PCQ, which acts as an interconnect to greatly magnify the NV-TLR coupling by several orders of magnitude, compared with the direct NV-TLR coupling (far below the linewidth of TLR with dozens of *kHz*) resulting from the vacuum fluctuations of the photons $2^{1,22}$ $2^{1,22}$ $2^{1,22}$. The TLR is made of a superconductor line interrupted by two capacitors at its ends. In the microwave domain, it can be treated as a quantum LC harmonic oscillator, $H^{TLR} = \omega_r (a^+ a + 1/2)$
(*b* - 1) where $\omega = \sqrt{1/I} C$ is the corresponding eigen- $(\hbar = 1)$, where $\omega_r = \sqrt{1/L_rC_r}$ is the corresponding eigen-
frequency with inductance *L*, and canacitance *C*. The PCO frequency with inductance *L^r* and capacitance *C^r* . The PCQ located at an antinode of TLR's magnetic field is formed by a superconducting loop interrupted by three Josephson junctions^{[23](#page-4-22)}. When the loop is biased by half a magnetic flux quantum, the device is an effective two-level qubit made up of two countercirculating persistent currents with the Hamiltonian $H^{PCQ} = \frac{\omega_0}{2} \sigma_z$. The NV can be modeled as a threefrom $H^2 = \frac{1}{2}\sigma_z$. The NV can be modeled as a three-
level system in the triplet ground-state subspace consisting of $\binom{3}{4}$, $m_s = 0$ and $\binom{3}{4}$, $m_s = \pm 1$. The Hamiltonian is $H^{NV} =$ $\gamma_e B_z S_z + D(S_z^2 - 2/3)$, where γ_e is the electronic gyromagnetic ratio $D/2\pi \sim 2.87 GHz$ is the zero-field splitting *R* is a netic ratio, $D/\sqrt{2\pi} \sim 2.87GHz$ is the zero-field splitting, B_z is a
perpendicular magnetic field at the NV and S, is the spin-1-z perpendicular magnetic field at the NV and *S ^z* is the spin-1-*z* operator.

The PCQ magnetically couples to TLR via mutual inductance, $H^{T-P} = -\hat{\mu} \cdot \hat{B}$, where $\hat{\mu}$ is the magnetic dipole of **PCO** induced by the persistent circulating currents and \hat{B} is PCQ induced by the persistent circulating currents and \hat{B} is the magnetic field at PCQ induced by the current in the central conductor of TLR. When $\omega_r \sim \omega_0$, we have $H^{T-P} =$
 $g(a^+ \tau^- + a \tau^+)$ after rotating wave approximation, where $g =$ $g(a^+\sigma^- + a\sigma^+)$ after rotating wave approximation, where $g = (I_{AB}r^2/d)\sqrt{(a+2I_{BA}r^2)}$ $r(L)$ is the radius (persistent circulating $(I_p\mu_0 r^2/d) \sqrt{\omega_r/2L_r}$, $r(I_p)$ is the radius (persistent circulating
current) of the PCO loop, and *d* is the distance between PCO current) of the PCQ loop, and *d* is the distance between PCQ and central conductor of TLR. The sizable changes of magnetic flux within the loop induced by I_p presented in the PCQ

lead to small shifts in the transition frequencies ($m_s \rightarrow \pm 1$) of $NV^{22,23}$ $NV^{22,23}$ $NV^{22,23}$ $NV^{22,23}$. Through this small change in magnetic field the PCQ can couple to the NV via Zeeman term, $H^{N-P} = \frac{1}{2} \eta \sigma_z S_z$, where $n = L \mu_0 \gamma_z / r$ where $\eta = I_p \mu_0 \gamma_e / r$.

The basic unit in our system is thus governed by the Hamil- $H_p^0 = H^{TLR} + H^{PCQ} + H^{NV} + H^{T-P} + H^{N-P}$. The photonic tunneling in our model can be realized by a central coupler^{[24](#page-4-23)} which serves as individual tunable quantum transducers to transfer photonic states between adjacent TLRs. We have presented a new paradigm for 2*D* TLR lattice coupled to solid-state spins. We have shown that specially engineered resonator lattice provides a practical platform to couple both individual spin and superconducting qubit, and engineer their interactions in a way that surpasses the limitations of current technologies. This can provide new insights to many-body physics.

III. MOTT-SUPERFLUID TRANSITION

We study the full Hamiltonian of the 2*D* square lattice by adding the on-site chemical potential and the nonlocal microwave photon hopping between adjacent sites. The Hamiltonian is given by

$$
H = \sum_{p} H_{p}^{0} + \sum_{\langle p,q \rangle} k_{\langle p,q \rangle} a_{p}^{+} a_{q} - \sum_{p} \mu_{p} N_{p}, \tag{1}
$$

where $k_{\langle p,q\rangle} = 2Z_0C_{\langle p,q\rangle}(\omega_r + \delta_p)(\omega_r + \delta_q)$ are photonic tunneling rates between resonators *p* and *q*, which are set by the tunable mutual capacitance $C_{\langle p,q\rangle}$ between resonator ends with characteristic impedance Z_0 and frequency shift δ_p due to ran-dom disorder^{[19](#page-4-18)}. Since $\omega_r \gg \delta_p$, one can assume that $k_{\langle p,q \rangle} =$ $k = 2Z_0C\omega_r^2$ without disorder for nearest-neighbor resonators,
and $k_{\text{max}} = 0$ for other resonator pairs, μ_r is the chemical poand $k_{\langle p,q\rangle} = 0$ for other resonator pairs. μ_p is the chemical potential at the *p*-th site. The conserved quantity in our system is the total number of excitations $N_p = a_p^+ a_p + \sigma_p^+ \sigma_p^- + \frac{1}{2} S_p^+ S_p^$ with $S^i(\sigma^i)$ the spin-1 (-1/2) operators $(i = x, y, \pm)$.
The photon-dependent eigenstates of the Hamilt

The photon-dependent eigenstates of the Hamiltonian H_p^0 is dressed states $|m_s, \pm, n\rangle$ with the eigenvalues $E_{|m_s, \pm, n\rangle} =$ $n\omega_r + (-\Delta + 2n\Delta \pm \sqrt{4ng^2 + \Delta^2 \pm 2\Delta\eta + \eta^2})/2 + \chi(m_s)$, where $\chi(m_s) = D(3 \times 1^{m_s} - 2) + m \propto R$ is the eigenenergy of H^{NV} $\chi(m_s) = D(3 \times 1^{m_s} - 2) + m_s \gamma_e B_z$ is the eigenenergy of H^{NV} .
Here $\Delta = (0, -\omega_0)$ is the defuning *n* is the number of ex-Here $\Delta = \omega_r - \omega_0$ is the detuning, *n* is the number of exclusions in the reconstant and $|\pm \rangle = (|\alpha \rangle + |\alpha \rangle)/\sqrt{2}$. In our citations in the resonator and $|\pm\rangle = (|e\rangle \pm |g\rangle)/\sqrt{2}$. In our case the dynamics is governed by the Jaynes-Cummings (IC) case, the dynamics is governed by the Jaynes-Cummings (JC) type of interaction, which enables the interconversion of qubit excitations and photons, and provides the effective on-site repulsion. Meanwhile, pairs of TLRs are coupled by the twosite Hubbard model via one-photon hopping. The difference between the Bose–Hubbard model $(BHM)²⁸$ $(BHM)²⁸$ $(BHM)²⁸$ and our model is that the conserved particles are the polaritons rather than the pure bosons in BHM, and the effective on-site repulsion $U_{\pm}(n) = E_{|m_s,\pm,n+1\rangle} - E_{|m_s,\pm,n\rangle}$ decreases with the growth of heater proposition number and $U_{\pm}(n) \rightarrow 0$ in the limit of large n and photon number, and $U_{+}(n) \rightarrow 0$ in the limit of large *n* and $\Delta = 0$, as shown in Fig. 2, while it is a constant in BHM.

The phase diagrams can be distinguished using the corresponding order parameters. Here we choose the SF order parameter $\psi = \langle a_p \rangle$ (set to be real) to differentiate between insulator-like and SF-like states. Using the mean-field

FIG. 3. (color online). The phase diagrams in $\mu \sim k$ plane for different sets of NV-PCQ coupling η and tunable magnetic fields B_z applied on NV, where μ and k are the chemical potential and the photon hopping rate. The common parameters are $g = 1$, $\omega_r = 200$, $D = 100$ and $\gamma_e = -10^3$. The other parameters are set as (a) $\eta = 0.01$,
 $B = 0.0005T$ and $\Delta = a$, the phase boundaries are plotted in (d) $B_7 = 0.0005T$ and $\Delta = g$, the phase boundaries are plotted in (d), where $\Delta = 2g$ (solid), *g* (dashed) and 0 (dot-dashed); (b) $\eta = 1.2$, $B_7 = -0.3$ and $\Delta = 0$, the phase boundaries are plotted in (e), where $B_z = -0.3T$ (solid), 0.0005*T* (dashed) and 0.3*T* (dot-dashed); (c) $\eta = 0.75$, $B_z = 0.3T$ and $\Delta = 0$, the phase boundaries are plotted in (f), where $\eta = 0.01$ (solid), $\eta = 0.75$ (dashed) and $\eta = 1.5$ (dotdashed).

theory^{[29](#page-5-1)} we decouple the hopping term as $a_p^+ a_q = \left\langle a_p^+ \right\rangle a_q + \left\langle a_p^+ \right\rangle a_q$ $a_p^+\left\langle a_q\right\rangle - \left\langle a_p^+\right\rangle \left\langle a_q\right\rangle$, the resulting mean-field Hamiltonian can then be written as a sum over single sites,

$$
H^{MF} = \sum_{p} [H_p^0 - zk\psi(a_p^+ + a_p) + zk\psi^2 - \mu_p N_p], \qquad (2)
$$

where $z = 4$ is the number of nearest neighbours. Noting that $[H^{MF}, S_z] = 0$, therefore, we can treat S_z in the mean-field
Hamiltonian as a c-number and $S_z = +1.0$ Minimizing the Hamiltonian as a *c*-number and $S_z = \pm 1, 0$. Minimizing the ground state energy of the Hamiltonian H^{MF} with respect to ψ for different values of *u* and *k* we obtain the mean field phase for different values of μ and k , we obtain the mean field phase diagram/boundary in the (μ, k) plane when η varies from the weak coupling regime $(\eta \ll g)$ to the strong coupling regime $(\eta > g)$ under the resonant/detuning case. The features of Fig. 3 are rich. The regime where $\psi = 0$ corresponds to the stable and incompressible MI lobes characterized by a fixed number of excitations at per site with no variance. In each MI lobe, due to the nonlinearity and anharmonicity in the spec-trum originating from the photon blockade effect^{[30](#page-5-2)}, the strong emitter-field interaction leads to an effective large polaritonpolariton repulsion which freezes out hopping and localizes polaritons at individual lattice sites. By contrast, strong hop-

FIG. 4. (color online). (a) The order parameter ψ in $\mu \sim k$ plane. The parameters are $g = 1$, $\omega_r = 200$, $D = 100$, $\gamma_e = -10^3$, $\eta = 0.01$, $B_z = 0.0005T$, $\alpha = 0.2$ and $\Delta = 0$. The corresponding phase boundaries 0.0005*T*, $\alpha = 0.2$, and $\Delta = 0$. The corresponding phase boundaries are plotted in (b), where the solid, dashed, and dot-dashed line denote $\alpha = 0, 0.2,$ and 0.24, respectively.

ping favours delocalization and condensation of the particles into the zero-momentum state, namely, $\psi \neq 0$ indicates a SF compressible phase with the stable ground state at each site corresponding to a coherent state of excitations.

Analogous to the BHM, the physical picture behind is that the MI-SF phase transition results from the interplay between polariton delocalization and on-site repulsive interaction. Therefore, the phase boundary primarily depends on the ratio of the photon-hopping rate to the on-site repulsion rate. When the on-site repulsion dominates over hopping, the system should be in the MI phase, otherwise the system will be in the SF phase. From the expression of the parameter η and *g*, we can find that reduction of the size of the PCQ loop will increase η but decrease *^g*, and the adjustment of the distance *^d* only affects TLR-PCQ interaction. Furthermore, the detuning ∆ is also tunable by varying the magnetic field applied on NV. In Fig. 3, one can find that the size of the MI lobes varies with ∆, with the largest Mott lobes found on resonance.

Further insight to the transition can be gained when the photon hopping with complex-valued amplitude exists in Eq. (1), where the hopping process becomes $-\sum_{p \mid p,q} k_{p}e^{i\phi_{p,q}}a_p^{\dagger}a_q$
with $\phi_{p,q} = -\phi_{p,q}$ and we set $k_{p,q} = k$. We emphasize that with $\phi_{\langle p,q\rangle} = -\phi_{\langle q,p\rangle}$ and we set $k_{\langle p,q\rangle} = k$. We emphasize that this process is possible if the intermediate coupling elements are used to break time-reversal symmetry $6,16,31$ $6,16,31$ $6,16,31$. The parameter $k_{(p,q)}e^{i\phi_{(p,q)}}$ provides a new regime in the dynamical evo-
lution of the system. The sum of the tunnaling phases along lution of the system. The sum of the tunneling phases along a closed loop surrounding the plaquette is $2(\phi_{p+1,q} + \phi_{p,q+1} - \phi_p)$ $\phi_{p,q} - \phi_{p+1,q+1} = 2\pi\alpha$, which is actually the flux quanta per plaquette. Assuming that αs are all equal, the total Hamiltonian under mean-field approximation reads

$$
H_{\alpha}^{MF} = \sum_{p} [H_p^0 - zk\psi \cos(2\pi\alpha)(a_p^+ + a_p)
$$

+ $zk\psi^2 \cos(2\pi\alpha) - \mu_p N_p].$ (3)

The results are exhibited in Fig. 4, we find that the boundary line gradually shifts to the right as α enhances in the interval $[0, 1/4]$. Because of the spatial variation of tunneling phase, the wave function of a polariton from one lattice site to another acquires a nontrivial phase (*Aharonov-Bohm* phase)^{[32](#page-5-4)}, which actually reduces the effective hopping rates.

FIG. 5. (color online). The order parameter ψ in $\mu \sim k$ plane under the different dissipative rates (a) $\Gamma = \kappa = 0.01$; (b) $\Gamma = \kappa = 0.05$; (c) $\Gamma = \kappa = 0.1$; (d) $\Gamma = \kappa = 0.15$. The other parameters are $g = 1$, $\omega_r = 200, D = 100, \gamma_e = -10^3, \eta = 0.01, B_z = 0.0005T, \text{ and } \Delta = 0.$

IV. DISSIPATIVE EFFECTS

Generally, nonequilibrium processes such as dissipative effect, are crucial in solid-state devices. We show that the signature of the localization-delocalization transition remains even in the presence of the engineered dissipation by the quantum trajectory method^{[33](#page-5-5)}. The non-Hermitian Hamiltonian is formulated as

$$
H_{de} = H^{MF} - \frac{i\Gamma}{2} \sum_{p} \sigma_p^+ \sigma_p^- - \frac{i\kappa}{2} \sum_{p} a_p^+ a_p, \tag{4}
$$

where $\kappa = 4Z_0^2 C_{out}^2 \omega_r^3$ is the decay rate of TLR, and Γ is the decay rate from the effective excited state $|e\rangle$ of PCO. In our where $\kappa = 4Z_0 C_{out} \omega_r$ is the decay rate of TLK, and T is the decay rate from the effective excited state $|e\rangle$ of PCQ. In our case, the dissipative effects result from the unavoidable interaction between the PCQ/TLR and the corresponding Markovian environment, for example, the interaction between the output of the TLR and the corresponding vacuum field will result in a photonic escape rate with κ to the continuum. Here the dissipative effects of NV are negligible, compared with κ and Γ. The phase diagrams under dissipative effects are displayed in Fig. 5. Once the hopping rate is increased beyond a critical value, the system is expected to undergo a nonequilibrium QPT from a MI phase, where the initial photon population is self-trapped, to a SF phase with the dynamical photon population imbalance coherently oscillating between pairs of TLRs^{[6](#page-4-15)}. Furthermore, another obvious feature is that the size of MI phase becomes larger as the growth of dissipative rates. Note that the effective nonlinearity at per site becomes stronger at lower exciton numbers, which implies that the dissipative effect (inducing the decrease of the exciton numbers) favours the MI phase. As a result, the dissipation results in the dynamical switching from SF phase to MI phase and causes the increment of the size of MI phase.

V. EXPERIMENTAL FEASIBILITY

Firstly, we briefly stress the relevant experimental progress. From theoretical standpoint, it is possible to fabricate large arrays to observe many-body physics of interacting polaritons since resonators and qubits can be made lithographically 34 . Actually, it is indeed feasible to couple over 200 (or 1000) TLRs with negligible disorders (on the order of a few parts in $10⁴$) in a 2*D* lattice using a 32*mm* × 32*mm* sample or a full two-inch wafer^{[6](#page-4-15)}. Secondly, how to probe quantum manybody states of light is still an open question in photonic quan-tum simulation^{[35](#page-5-7)}. The previous works^{[2](#page-4-4)[,36](#page-5-8)} suggested to measure the individual TLR through mapping the excitations into the qubit followed by state-selective resonance fluorescence spectrum, but a remaining technical challenge is the realization of high-efficiency photon detectors. Alternatively, the local statistical property of TLR can be analyzed readily using combined techniques of photon-number-dependent qubit transition $37,38$ $37,38$ and fast readout of the qubit state through a sep-arate low-Q resonator mode^{[39](#page-5-11)}, for which the high-efficiency photon detectors are not required. Experimentally, transmission and reflection measurements for circuit QED arrays have been implemented successfully in small system with one or two resonators^{[7,](#page-4-24)[38](#page-5-10)}. Therefore, in order to distinguish between different phases of the system, one can also experimentally probe beyond transmission, such as two-tone spectroscopy and second-order coherence function (photon statistics) to reveal additional information. The tunability of coupling strengths in our system enables one to measure these quantities relatively straightforwardly.

Finally, we survey the relevant experimental parameters. Given the flexibility of circuit QED, we can access a wide range of tunable experimental parameters for TLR-PCQ coupling strength *g* and hopping rate $k_{(p,q)}$. Taking $L_r = 2 nH$, $ω_r/2π = 6 GHz, I_p = 600 nA, and r = 0.2 μm, we get$ $\eta/2\pi \simeq 140 \; KHz$ and $g/2\pi \sim [1.8, 180] \; MHz$ when the distance d varies from 5 μ *m* to 50 *nm*. Furthermore, the hopping rate $k_{\langle p,q \rangle}$ depends on the tunable mutual capacitance $C_{\langle p,q \rangle}$ between resonator ends. In Ref^{19} Ref^{19} Ref^{19} , the authors measured devices with photon hopping rates $k/2\pi$ form 0.8 MHz to 31 MHz in resonators lattices. On the other hand, the electron-spin relaxation time T_1 of NV ranges from 6 ms at room temperature^{[41](#page-5-12)} to 28 ∼ 265 s at low temperature^{[42](#page-5-13)}. In addition, later experi-mental progress^{[43](#page-5-14)} with isotopically pure diamond has demonstrated a longer dephasing time to be $T_2 = 1.8$ ms. Therefore, the dissipation and decoherence of NV are negligible.

VI. CONCLUSION

We have devised a concrete hybrid system to engineer photonic MI-SF phase transition in a 2*D* square lattice of TLRs coupled to a single NV encircled by a PCQ. We find that the interplay between the on-site repulsion and the nonlocal tunneling leads to the photonic localization-delocalization transition. In the presence of dissipation, the phase boundary can be obtained by the mean-field approach and the quantum jump technique. Facilitated by high levels of connectivity in circuit QED, experiments combining both scalability and long coherence times are expected in the coming few years, at that stage the investigation of photonic QPT using TLR lattice systems can therefore be easier to realize.

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