Exploring Unconventional Hubbard Models with Doubly Modulated Lattice Gases

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Recent experiments show that periodic modulations of cold atoms in optical lattices may be used to engineer and explore interesting models. We show that double modulation combining lattice shaking and modulated interactions allows for the engineering of a much broader class of lattice with correlated hopping, which we study for the particular case of one-dimensional systems. We show, in particular, that by using this double modulation it is possible to study Hubbard models with asymmetric hopping, which, contrary to the standard Hubbard model, present insulating phases with both parity and string order. Moreover, double modulation allows for the simulation of lattice models in unconventional parameter regimes, as we illustrate for the case of the spin-1/2 Fermi-Hubbard model with correlated hopping, a relevant model for cuprate superconductors.

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Introduction.—Ultracold gases in optical lattices have attracted a lot of attention as emulators of fundamental models of quantum many-body systems. Their unprecedented levels of controllability, tunability, and cleanness have permitted the realization of Hubbard models with cold atoms [1–3], the creation of synthetic magnetic fields in neutral lattice gases [4,5], first steps towards the emulation of quantum magnetism [6–12], and more [13,14]. Recent progress in measurement techniques has allowed for single-site resolved detection [15,16], which has permitted a deeper insight into the properties of Mott insulators (MI) [16–18] and the dynamical properties of cold lattice gases [19].

The possibility of tuning parameters in real time in cold lattice gases has aroused particular interest. Fast periodic modulations provide a new tool for the engineering of relevant lattice models [4,5,20–29]. In particular, a fast-enough modulation of the lattice position (lattice shaking) results in the effective change of the tunneling rate [20] allowing, for example, driving the superfluid (SF) to MI transition [24], inducing photon-assisted hopping in tilted lattices [22], simulating frustrated classical magnetism [25], generating gauge potentials [28], and inducing effective ferromagnetic domains [29]. Moreover, a fast modulation of the interparticle interactions results in an effective hopping that depends on the occupation differences at neighboring sites [30–33] and may induce density-dependent gauge fields [34].

In this Letter, we show that double modulation (DM), i.e., the combination of lattice shaking and periodically modulated interactions, permits the selective control of different hopping processes, hence, allowing for the engineering of a broad range of lattice models that cannot be realized with either lattice shaking or modulated interaction

alone. DM permits, in particular, the generation of mirror-asymmetric tunneling, which may result in insulators with both parity and string order. Moreover, DM permits the exploration of parameter regimes which cannot be reached with a single modulation, as we illustrate for the relevant case of spin-1/2 lattice fermions with correlated hopping.

Double modulation.—We consider a lattice gas with periodic DM. Whereas lattice shaking results from the displacement of the lattice [20], periodic interactions may be induced by modulating an externally applied magnetic field in the vicinity of a Feshbach resonance (see Refs. [30,32,33] for details). We focus below on 1D lattices, although the engineering possibilities of DM may be extended to higher dimensions as well, and in 2D lattices, elliptic shaking [35] may be employed to induce even richer lattice models. In the experimentally relevant scenario in which only the lowest Bloch band is relevant [36], Bose gases are described in the lattice reference frame by the time-dependent Bose-Hubbard model (BHM):

$$\hat{H} = -J \sum_{\langle i,j \rangle} \hat{b}_{i}^{\dagger} \hat{b}_{j} + \frac{U(t)}{2} \sum_{j} \hat{n}_{j} (\hat{n}_{j} - 1) + F(t) \sum_{j} j \hat{n}_{j}, \quad (1)$$

where \hat{b}_i (\hat{b}_i^{\dagger}) annihilates (creates) a boson at site i, $\hat{n}_i = \hat{b}_i^{\dagger} \hat{b}_i$, J is the tunneling parameter, U(t) is the time-dependent interaction strength, and F(t) is a time-dependent tilting amplitude resulting from the lattice shaking. Both the interaction and the tilting term are periodically modulated, $U(t) = U_0 + U_1 f_U(t) = U_0 + U_1 f_U(t+T)$ and $F(t) = F_1 f_F(t) = F_1 f_F(t+T)$. Both $f_{U,F}(t)$ are unbiased, i.e., $\int_t^{t+T} f_{U,F}(\tau) d\tau = 0$. If $\omega = 2\pi/T \gg U_0/\hbar$, J/\hbar , Floquet analysis may be employed to integrate the modulations, obtaining an effective time-independent Hamiltonian (see, e.g., Refs. [20,32] or

Ref. [33] for an equivalent derivation for the Fermi-Hubbard model):

$$\hat{H}_{\text{eff}} = -J \sum_{\langle i,j \rangle} \hat{b}_i^{\dagger} \mathcal{F}(i-j, \hat{n}_i - \hat{n}_j) \hat{b}_j + \frac{U_0}{2} \sum_j \hat{n}_j (\hat{n}_j - 1),$$
(2)

where $\mathcal{F}=(1/T)\int_0^T e^{i/\hbar}\int_0^t [F_1f_F(\tau)(i-j)+U_1f_U(\tau)(\hat{n}_i-\hat{n}_j)]d\tau dt$. We illustrate below the engineering possibilities allowed by DM for the particular case of $f_U(t)=f_F(t)=\cos(\omega t)$. Note, however, that different frequencies and/or functional forms for the two modulations may allow for an even more versatile engineering. The above-mentioned choice implies $\mathcal{F}=\mathcal{J}_0((F_1/\hbar\omega)(i-j)+(U_1/\hbar\omega)(\hat{n}_i-\hat{n}_j))$ in Eq. (2), where \mathcal{J}_n is the nth-order Bessel function [37].

Hopping channels.—We denote as $J_{(n_i,n_j)\leftrightarrow(n_i+1,n_j-1)}$ the hopping rate from site j with n_j particles before the hop to site i with initially n_i particles. In 1D lattices, a hop to the left is characterized by the rate

$$\frac{J_{(n_j,n_{j+1})\leftrightarrow(n_j+1,n_{j+1}-1)}}{J} = \mathcal{J}_0\left(\frac{U_1}{\hbar\omega}(n_j-n_{j+1}+1) - \frac{F_1}{\hbar\omega}\right),$$

whereas a hop to the right is given by

$$\frac{J_{(n_j,n_{j+1})\leftrightarrow(n_j-1,n_{j+1}+1)}}{J} = \mathcal{J}_0\left(\frac{U_1}{\hbar\omega}(n_{j+1}-n_j+1) + \frac{F_1}{\hbar\omega}\right).$$

Note that for sole lattice shaking $(U_1 = 0)$, hops are mirror symmetric since \mathcal{J}_0 is even. The same is true for solely modulating interactions $(F_1 = 0)$. DM allows for breaking mirror symmetry. Especially relevant at low fillings are $J_{(0,1)\leftrightarrow(1,0)}=J_{(1,2)\leftrightarrow(2,1)}=J\mathcal{J}_0(F_1/\hbar\omega),$ $J_{(1,1)\leftrightarrow(2,0)} = J\mathcal{J}_0((U_1 - F_1)/\hbar\omega), \text{ and } J_{(1,1)\leftrightarrow(0,2)} =$ $J\mathcal{J}_0((U_1+F_1)/\hbar\omega)$. One may observe two important features: (i) contrary to the standard BHM $(F_1 = U_1 = 0)$, in general, $J_{(1,1)\leftrightarrow(2,0)} \neq J_{(1,1)\leftrightarrow(0,2)}$, and (ii) $J_{(1,1)\leftrightarrow(2,0)}/$ $J_{(0,1)\leftrightarrow(1,0)}$ and/or $J_{(1,1)\leftrightarrow(0,2)}/J_{(0,1)\leftrightarrow(1,0)}$ may be larger than 1, an impossibility for $U_1 = 0$ and/or $F_1 = 0$. These peculiar features have crucial consequences. As shown below, (i) DM may result in insulators with both parity and string order, whereas (ii) it also allows for the study of a much richer phase diagram for lattice gases, compared to the case of either shaking or interaction modulation.

Insulators with finite parity and string order.—In the standard BHM, the MI with unit occupation is characterized by doublon-holon pairs in a sea of singly occupied sites. The pairs result from the $J_{(1,1)\leftrightarrow(2,0)}=J_{(1,1)\leftrightarrow(0,2)}$ hops. The 1D MI at unit filling presents nonlocal parity order $\mathcal{O}_P^2 \equiv \lim_{|i-j|\to\infty} \langle (-1)^{\sum_{i< l < j} \delta \hat{n}_i} \rangle > 0$, with $\delta \hat{n}_j = 1 - \hat{n}_j$ [38], due to the appearance of doublon and holon defects in pairs, as recently revealed experimentally [18].

Another important nonlocal order in 1D is string order, $\mathcal{O}_S^2 \equiv \lim_{|i-j| \to \infty} - \langle \delta \hat{n}_i (-1)^{\sum_{i < l < j} \delta \hat{n}_l} \delta \hat{n}_j \rangle$ [39]. A nonvanishing \mathcal{O}_S^2 characterizes the Haldane insulator (HI) predicted in polar lattice gases [39] and bosons in frustrated lattices [40]. In the HI, the position of the defects and the separation between them is not fixed, but starting with a doublon, the next defect along the chain is a holon, the next a doublon, and so on. This diluted "antiferromagnetic" order is characterized by $\mathcal{O}_S^2 > 0$. However, in a HI, defects are not paired, and, hence, $\mathcal{O}_P^2 = 0$. Conversely, in the standard BHM the MI presents $\mathcal{O}_S^2 = 0$ due to the equal probability of having 20 and 02 pairs.

DM allows for the selective reduction or cancellation of $J_{(1,1)\leftrightarrow(2,0)}$ [or, alternatively, $J_{(1,1)\leftrightarrow(0,2)}$]. As a result, inside the MI, pairs may be produced dominantly, or even solely, following a 02 (or 20) order. The system may then present both $\mathcal{O}_P^2 > 0$, due to the pairwise creation of defects, and $\mathcal{O}_S^2 > 0$, due to the dominantly antiferromagnetic order of the defects. This is best understood from the case in which all hoppings vanish except $J_{(1,1)\leftrightarrow(0,2)}$ [41]. For unit filling, since the defect-free MI at J=0 just develops pairs 02 at finite J, we map into an effective spin-1/2 system, with $|0,2\rangle_{j,j+1} \equiv |\uparrow\rangle_j$ and $|1,1\rangle_{j,j+1} \equiv |\downarrow\rangle_j$, obtaining an effective spin model:

$$H_{S} = \sum_{j} J_{(1,1)\leftrightarrow(0,2)}(\hat{S}_{j}^{+} + \hat{S}_{j}^{-}) + U_{0}\left(\frac{1}{2} + \hat{S}_{j}^{z}\right) + \Delta\left(\frac{1}{2} + \hat{S}_{j}^{z}\right)\left(\frac{1}{2} + \hat{S}_{j+1}^{z}\right), \tag{3}$$

where $\hat{S}_{i}^{\pm,z}$ are spin-1/2 operators, and we add a spin-spin interaction $\Delta \to \infty$ to project out neighboring up spins, which have no physical meaning. Interestingly, Hamiltonian (3) corresponds to an Ising model with transverse and longitudinal magnetic fields in the vicinity of its tricritical point between, ferro-, antiferro-, and paramagnetic phases [42]. For a chain of length L and periodic boundary conditions, the ground state of (3) is a linear combination of states with m spin-ups, i.e., m defect pairs, of the form $\sum_{m=0}^{L/2} c_m |m\rangle$, where $|m\rangle \propto (PS^+)^m |0\rangle$ with Pthe projector excluding states with two neighboring up spins, $S^+ = \sum_i S_i^+$, and $|0\rangle$ the state with all spins down (i.e., a defect-free MI). String and parity order may then be expressed as statistical moments of the number of defect pairs: $\mathcal{O}_P^2 \approx \langle (2m-1)^2 \rangle$ and $\mathcal{O}_S^2 \approx \langle 2m^2(m-2) \rangle$. Both orders may, hence, coexist. Note that with an increasing number of pairs, i.e., decreasing ratio $U_0/J_{(1,1)\leftrightarrow(0,2)}$, \mathcal{O}_S^2 increases whereas \mathcal{O}_P^2 decreases, as observed in our simulations of Eq. (2).

Figure 1 (top) illustrates the rich physics that results from the selective control of the different hopping rates allowed by DM. For a fixed $U_0/J=1.2$ and $U_1/\hbar\omega=3$, we

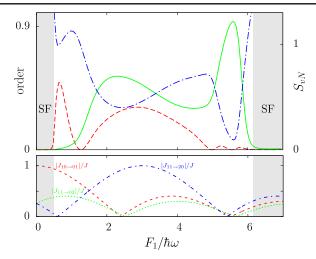


FIG. 1 (color online). (top) Infinite DMRG [43] results for the effective model (2) (with four bosons per site, keeping up to M=200 states) for \mathcal{O}_P^2 (solid), $4\mathcal{O}_S^2$ (dashed), and S_{vN} (dot dashed) as a function of $F_1/\hbar\omega$ for $U_0/J=1.2$, unit filling, and $U_1/\hbar\omega=3$. In the shaded SF regions, we find a logarithmic divergence of S_{vN} with the number of kept matrix states M. (bottom) Relevant hopping rates.

analyze, by means of density matrix renormalization group (DMRG) [43] calculations, the different order parameters for a varying $F_1/\hbar\omega$. The relevant tunneling rates are depicted in Fig. 1 (bottom). One observes both that the hopping is generally asymmetric, $J_{(1,1)\leftrightarrow(2,0)} \neq J_{(1,1)\leftrightarrow(0,2)}$, and the existence of parameter regimes at which $J_{(0,1)\leftrightarrow(1,0)} < J_{(1,1)\leftrightarrow(2,0)}$ and/or $J_{(1,1)\leftrightarrow(0,2)}$. A first consequence of the variation of the hopping rates with F_1 is clearly the appearance of SF to insulator transitions (at $F_1/\hbar\omega \simeq 0.5$ and 6.2). In the SF regime, the excitation gap vanishes, the entanglement entropy S_{vN} [44] shows a logarithmic divergence with the system size (not shown here), and correlations decay algebraically as expected for a Luttinger liquid. Within the insulator regions, the entanglement entropy is finite and correlations decay exponentially. As expected from the discussion above, insulator phases with $\mathcal{O}_P^2 > 0$ and $\mathcal{O}_S^2 > 0$ occur due to the hopping asymmetry. Note that \mathcal{O}_S^2 increases when the hopping asymmetry grows, disappearing at those F_1 values at which $J_{(1,1)\leftrightarrow(2,0)} = J_{(1,1)\leftrightarrow(0,2)}$ (at $F_1/\hbar\omega \approx 1.2, 5.2$, and 5.5). When $J_{(1,1)\leftrightarrow(0,2)}, J_{(1,0)\leftrightarrow(0,1)} \ll J_{(1,1)\leftrightarrow(2,0)}$, we recover the extreme case discussed above, and \mathcal{O}_P^2 and \mathcal{O}_S^2 are large (e.g., at $F_1/\hbar\omega \approx 2.4$). Interestingly, within the insulator regions, when the hopping is asymmetric and $J_{(1,1)\leftrightarrow(2,0)}$, $J_{(1,1)\leftrightarrow(0,2)} \ll J_{(1,0)\leftrightarrow(0,1)}$, we obtain $\mathcal{O}_P^2 \ll \mathcal{O}_S^2$. This parameter region is characterized by, to a large extent, broken defect pairs but still antiferromagnetic defect order; i.e., in this region the insulator rather behaves as a HI. We find, however, no gapless region that would mark a transition between a HI (with low \mathcal{O}_P^2) and a MI (with finite \mathcal{O}_S^2) [38].

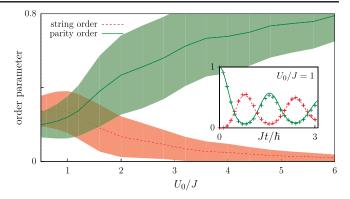


FIG. 2 (color online). Dynamics of \mathcal{O}_P^2 (solid) and \mathcal{O}_S^2 (dashed) after a sudden onset of DM with $F_1/\hbar\omega=2.4$, $U_1/\hbar\omega=3$ for a system initially prepared as a deep Mott insulator. The inset shows the dynamics for $U_0/J=1$. The curves are obtained using the effective model (2), whereas the crosses indicate the results directly obtained from Eq. (1). In the main figure we depict the time average of the orders for $2 < Jt/\hbar < 6$ as a function of U_0 for the same F_1 and U_1 . The shaded regions indicate the variances of the orders associated to the dynamics after the quench. In the DMRG simulations of Eq. (1), we employ L=60 sites, N=60 particles, and $\hbar\omega=20\pi J$; the simulations of the effective model (2) are performed in an infinite scheme [43]. In both cases we keep up to M=400 matrix states.

Dynamics after switching the double modulation.—Up to this point, we have analyzed the ground-state properties of the time-independent model (2). It is, however, interesting and experimentally relevant to investigate the dynamics following the switching of DM. For simplicity, we consider at times t < 0 a large U_0/J at unit filling in absence of any modulation, such that the system is in a defect-free MI $(\mathcal{O}_P^2 = 1, \mathcal{O}_S^2 = 0)$. At t = 0, U_0 is set to a final value, and the sinusoidal DM is abruptly switched on (to values $U_1/\hbar\omega = 3$ and $F_1/\hbar\omega = 2.4$ in Fig. 2), which results in $J_{(0,1)\leftrightarrow(1,0)}, J_{(1,1)\leftrightarrow(0,2)} \ll J_{(1,1)\leftrightarrow(2,0)}$. We have performed DMRG simulations of the dynamics employing both the original time-dependent Hamiltonian (1) and the effective one (2). As depicted in the inset of Fig. 2, both models provide identical results, showing the validity of the effective model for describing the dynamics. After the quench, the creation of defects abruptly reduces \mathcal{O}_P^2 and increases \mathcal{O}_S^2 and then subsequently develops an oscillatory dynamics. Figure 2 shows the time average of \mathcal{O}_P^2 and \mathcal{O}_S^2 as a function of U_0 (the shadowed regions indicate the variance of the orders related to the oscillations characterizing the real time dynamics). The average values present a similar qualitative dependence as that expected from the stationary state, with growing \mathcal{O}_P^2 and decreasing \mathcal{O}_S^2 for larger U_0/J (see Fig. 3 of Ref. [18] for an analysis of the growth of \mathcal{O}_P^2 for increasing U_0/J without periodic modulations). These results, hence, show that even an abrupt start of the double modulation transforms an initial defect-free MI into an insulator with nonvanishing timeaveraged \mathcal{O}_P^2 and \mathcal{O}_S^2 .

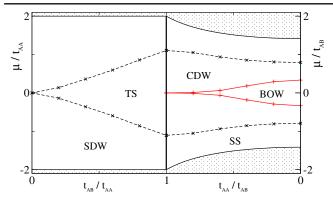


FIG. 3 (color online). Phase diagram for the Fermi-Hubbard model as a function of μ/t_{AA} and t_{AB}/t_{AA} for $U_0=0$ (see text). The dashed line denotes values of μ for which the Luttinger-liquid parameter $K_{\rho}=1$ extracted from the long wavelength behavior of the static charge structure factor [53] for $t_{AA} \neq t_{AB}$. This is consistent with the perturbative result from Ref. [48] for $t_{AA}=t_{AB}$. The size for the BOW phase corresponds to the charge gap at unit filling. The shaded regions denote the vacuum or the fully occupied state. All results are extrapolated to the thermodynamic limit from open boundary DMRG calculations with up to L=144 sites.

Exploring the complete phase diagram of fermions with correlated hopping.—As mentioned above, double modulation allows for $J_{(1,1)\leftrightarrow(2,0)}/J_{(0,1)\leftrightarrow(1,0)}$ and/or $J_{(1,1)\leftrightarrow(0,2)}/J_{(0,1)\leftrightarrow(1,0)}$ to be larger than 1, and, hence, for exploring novel quantum phases unreachable with $U_1=0$ and/or $F_1=0$. This is best illustrated by the spin-1/2 Fermi-Hubbard model with lattice shaking and modulated interactions $[U_0+U_1\cos{(\omega t)}]\sum_j \hat{n}_{j\uparrow}\hat{n}_{j\downarrow}$. Considering for simplicity a spin-independent lattice and mirror-symmetric hopping, $\mathcal{J}_0[(F_1+U_1)/\hbar\omega]=\mathcal{J}_0[(F_1-U_1)/\hbar\omega]$, we reach for sufficiently fast modulations the effective Hamiltonian:

$$H_{F} = U_{0} \sum_{j} \hat{n}_{j\uparrow} \hat{n}_{j\downarrow} - \sum_{\langle i,j\rangle,\sigma=\uparrow,\downarrow} (\hat{c}^{\dagger}_{i\bar{\sigma}} \hat{c}_{j\bar{\sigma}} + \text{H.c.}) \mathcal{P}(\hat{n}_{i\sigma}, \hat{n}_{j\sigma}),$$

$$(4)$$

with $\bar{\sigma}=-\sigma$, and $\mathcal{P}(\hat{n}_{i\sigma},\hat{n}_{j\sigma})\equiv t_{AA}(1-\hat{n}_{i\sigma})(1-\hat{n}_{j\sigma})+t_{BB}\hat{n}_{i\sigma}\hat{n}_{j\sigma}+t_{AB}[\hat{n}_{i\sigma}(1-\hat{n}_{j\sigma})+\hat{n}_{j\sigma}(1-\hat{n}_{i\sigma})]$, where $\hat{c}_{j\sigma}$ annihilates a fermion of spin σ at site j, $\hat{n}_{j\sigma}=\hat{c}_{j\sigma}^{\dagger}\hat{c}_{j\sigma}$, $t_{AA}=t_{BB}=J\mathcal{J}(F_1/\hbar\omega)$, and $t_{AB}=J\mathcal{J}_0((U_1+F_1)/\hbar\omega)$. Hamiltonian (4) has been extensively studied as a model for cuprate superconductors [45–49]. Figure 3 shows the grand-canonical phase diagram for $U_0=0$ [50]. For $t_{AB}< t_{AA}$, two spin-massless phases occur [48]: one with dominant triplet superconducting (TS) [51] correlations $\langle \mathcal{Q}_{0+}^{\dagger}\mathcal{Q}_{j+}\rangle$, with $\mathcal{Q}_{j\pm}\equiv c_{j+1\downarrow}c_{j\uparrow}\pm c_{j+1\uparrow}c_{\downarrow}$ and another with dominant spin-density wave (SDW) correlations, $(-1)^j\langle\hat{n}_{0-}\hat{n}_{j-}\rangle$, with $\hat{n}_{j\pm}\equiv\hat{n}_{j\uparrow}\pm\hat{n}_{j\downarrow}$. At $t_{AB}=0$, the TS phase has vanishing Drude weight (Kohn metal

[45,46,52]), and the SDW at $\bar{n} \equiv \langle \hat{n}_{j+} \rangle > 1 (<1)$ is a metal without holons (doublons) [45].

Sole modulation of interactions only permits $0 \le t_{AB}/t_{AA} \le 1$ [33]. DM allows for $t_{AB}/t_{AA} > 1$, for which three spin-massive phases occur [48]: a gapped phase (both in spin and density sectors) at $\bar{n}=1$ with bond-ordering wave (BOW) order, $\langle \mathcal{B}_0^{\dagger} \mathcal{B}_j \rangle$ with $\mathcal{B}_j \equiv \sum_{\sigma} (\hat{c}_{j+1\sigma}^{\dagger} \hat{c}_{j\sigma} + \text{H.c.})$, a gapless phase with dominant density wave correlations, $(-1)^j \langle \hat{n}_{0+} \hat{n}_{j+} \rangle$, and a gapless phase with dominant singlet-superconducting correlations, $\langle \mathcal{Q}_{0-}^{\dagger} \mathcal{Q}_{j-} \rangle$. Interestingly, the gapped BOW phase occurs even for $U_0=0$ due to the effective repulsion induced by the density-dependent hopping.

Outlook.—DM of cold lattice gases allows for the precise control of selected hopping processes. Such a control permits the realization of quantum phases with unconventional properties. In particular, mirror-asymmetric hopping results in insulating 1D phases with both parity and string orders, which may be revealed using *in situ* site resolved imaging [54]. We have also shown that DM may be used to simulate lattice models in otherwise unreachable regimes, as shown for the relevant case of the spin-1/2 Fermi-Hubbard model with correlated hopping.

We have considered, for simplicity, homogeneous systems. With an overall confinement V(j) that varies slowly enough from site to site, the grand-canonical phase diagram maps into a spatial distribution through the local chemical potential $\mu_i = \mu - V(j)$. In particular, the BOW phase results in a density plateau that may be revealed using single-site resolution [16,17]. Moreover, whereas for $U_0 = 0$, the spin gap opens at $t_{AB,cr} = t_{AA}$ at any \bar{n} , for $U_0 > 0$, $t_{AB,cr}$ depends on U_0 and \bar{n} (for low U_0/t_{AA} , $t_{AB,cr} \simeq U_0/[8(\bar{n}-1)\cos(\pi\bar{n}/2) - 16\sin(\pi\bar{n}/2)/\pi]$ [48]). As a result, for fixed t_{AB} , t_{AA} , and U_0 , the spin gap opens at a critical μ_{cr} . For an overall confinement, there is, hence, a spatial boundary between spin-gapped and spin-gapless phases, which may be revealed by creating spin excitations in the gapless region using a spin-dependent potential and observing the reflection of the excitations at the boundary [55].

In this Letter, we have focused on the simplest double modulation possible in which both fields were oscillating in phase. A natural and nontrivial extension of this work would be in the analysis of the physics emerging when the two modulations are different and time-reversal symmetry is broken. Another important extension of this work would be in the application of DM to 2D and 3D systems, where the combination of elliptical lattice shaking [28] and periodically modulated interactions may lead to an even richer physics.

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Leibniz Universität Hannover. Computer simulations for the inset of Fig. 2, run on the cluster Perseus at the Department of Theoretical Physics of the University of Geneva, employed the Algorithms and Libraries for Physics Simulations libraries [56,57] and the code IVAN.

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