

Quantum simulation of field theories with ultracold atoms

A. Trombettoni
(University of Trieste & INFN)

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Outline

- **Ultracold atoms as quantum simulators for field theories**
- **Symmetry-locked superfluid phases**
- **Link models**

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- **Symmetry-locked superfluid phases**
- **Link models**

The ultracold family: Bosons & Fermions

																		Key																				
																		11	Atomic number																			
																		Na	Element symbol																			
																		Sodium	Element name																			
																		Average atomic mass*																				
1	1A	1	2															13	14	15	16	17	18															
1		H																5	6	7	8	9	10															
		Hydrogen																Boron	Carbon	Nitrogen	Oxygen	Fluorine	Neon															
		1.01																10.81	12.01	14.01	16.00	19.00	20.18															
2	2A	3	4																																			
		Li	Be																																			
		Lithium	Beryllium																																			
		6.94	9.01																																			
3		11	12																																			
		Na	Mg																																			
		Sodium	Magnesium																																			
		22.99	24.31																																			
4		19	20	21	22	23	24	25	26	27	28	29	30	31	32	33	34	35	36																			
		K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr																			
		Potassium	Calcium	Scandium	Titanium	Vanadium	Chromium	Manganese	Iron	Cobalt	Nickel	Copper	Zinc	Gallium	Germanium	Arsenic	Selenium	Bromine	Krypton																			
		39.10	40.08	44.96	47.87	50.94	52.00	54.94	55.85	58.93	58.69	63.55	65.39	69.72	72.61	74.92	78.96	79.90	83.80																			
5		37	38	39	40	41	42	43	44	45	46	47	48	49	50	51	52	53	54																			
		Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe																			
		Rubidium	Strontium	Yttrium	Zirconium	Niobium	Molybdenum	Technetium (98)	Ruthenium	Rhodium	Palladium	Silver	Cadmium	Indium	Tin	Antimony	Tellurium	Iodine	Xenon																			
		85.47	87.62	88.91	91.22	92.91	95.94	(98)	101.07	102.91	106.42	107.87	112.41	114.82	118.71	121.76	127.60	126.90	131.29																			
6		55	56	57	72	73	74	75	76	77	78	79	80	81	82	83	84	85	86																			
		Cs	Ba	La	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn																			
		Cesium	Barium	Lanthanum	Hafnium	Tantalum	Tungsten	Rhenium	Osmium	Iridium	Platinum	Gold	Mercury	Thallium	Lead	Bismuth	Polonium (209)	Astatine (210)	Radon (222)																			
		132.91	137.33	138.91	178.49	180.95	183.84	186.21	190.23	192.22	195.08	196.97	200.59	204.38	207.2	208.98	(209)	(210)	(222)																			
7		87	88	89	104	105	106	107	108	109																												
		Fr	Ra	Ac	Rf	Db	Sg	Bh	Hs	Mt																												
		Francium (223)	Radium (226)	Actinium (227)	Rutherfordium (261)	Dubnium (262)	Seaborgium (266)	Bohrium (264)	Hassium (269)	Mendelevium (268)																												
																		58	59	60	61	62	63	64	65	66	67	68	69	70	71							
																		Ce	Pr	Nd	Pm	Sm	Eu	Gd	Tb	Dy	Ho	Er	Tm	Yb	Lu							
																		Cerium	Praseodymium	Neodymium	Promethium (145)	Samarium	Europium	Gadolinium	Terbium	Dysprosium	Holmium	Erbium	Thulium	Ytterbium	Lutetium							
																		140.12	140.91	144.24	(145)	150.36	151.96	157.25	158.93	162.50	164.93	167.26	168.93	173.04	174.97							
																		90	91	92	93	94	95	96	97	98	99	100	101	102	103							
																		Th	Pa	U	Np	Pu	Am	Cm	Bk	Cf	Es	Fm	Md	No	Lr							
																		Thorium	Protactinium	Uranium	Neptunium (237)	Plutonium (244)	Americium (243)	Curium (247)	Berkelium (247)	Californium (251)	Einsteinium (252)	Fermium (257)	Mendelevium (258)	Nobelium (259)	Lawrencium (262)							
																		232.04	231.04	238.03	(237)	(244)	(243)	(247)	(247)	(251)	(252)	(257)	(258)	(259)	(262)							

* If this number is in parentheses, then it refers to the atomic mass of the most stable isotope.

Typical values: $\left\{ \begin{array}{l} \text{numbers: } 10^3 - 10^6 \text{ atoms} \\ \text{temperatures: } 10 - 100 \text{ nanoKelvin} \\ \text{sizes: } 1 - 50 \text{ micrometers} \end{array} \right.$

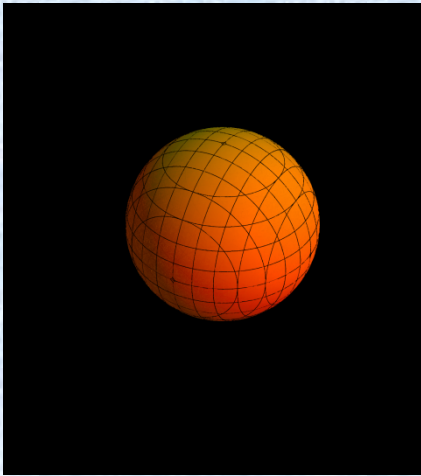
Controlling the system...

- **Bosons and/or fermions**
- **Geometry (1D / 2D)**
- **Long-range interactions**
- **Add disorder**
- **Time-dependence (and to a certain extent space-dependence) of the parameters of the Hamiltonian**
- **Simulate a magnetic field through a rotation or with optical tools**
- **Explicit tuning of the interactions via Feshbach resonances**
- **Optical lattices (i.e., periodic potentials and minima of the potential located on a lattice)**

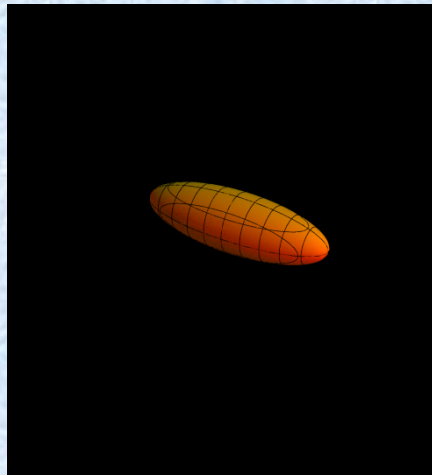
Experimental setup (I)

Magnetic harmonic potential: $V(x, y, z) = \frac{1}{2} m (\omega_x^2 x^2 + \omega_y^2 y^2 + \omega_z^2 z^2)$

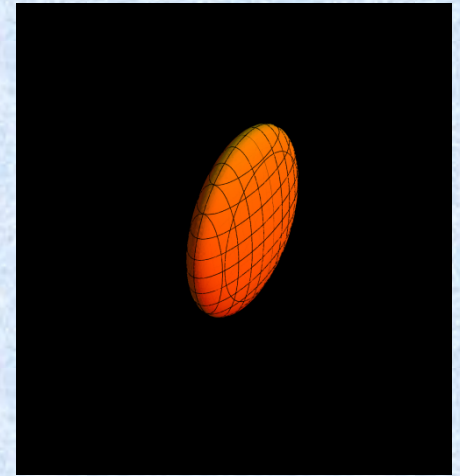
$$\omega_x = \omega_y = \omega_z$$



$$\omega_x \ll \omega_y = \omega_z$$



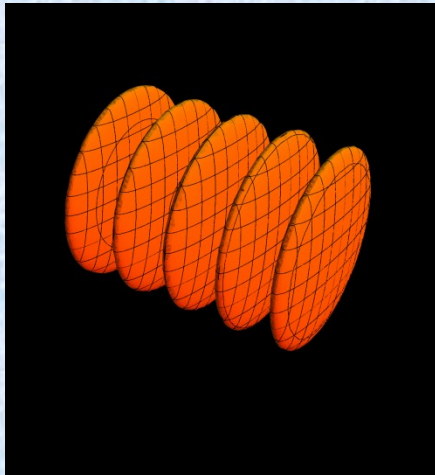
$$\omega_x \gg \omega_y = \omega_z$$



Experimental setup (II)

Using optical lattices:

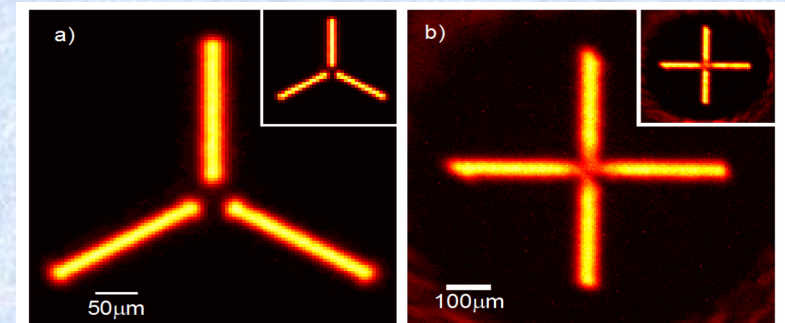
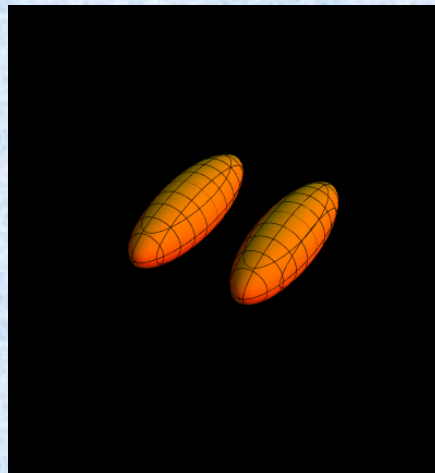
$$V = V_0 \cos^2(kx)$$



$$V = V_0 \left[\cos^2(kx) + \cos^2(ky) \right]$$



Other configurations (like ladders, coupled cigars or stars) are as well possible:



Quantum Simulations

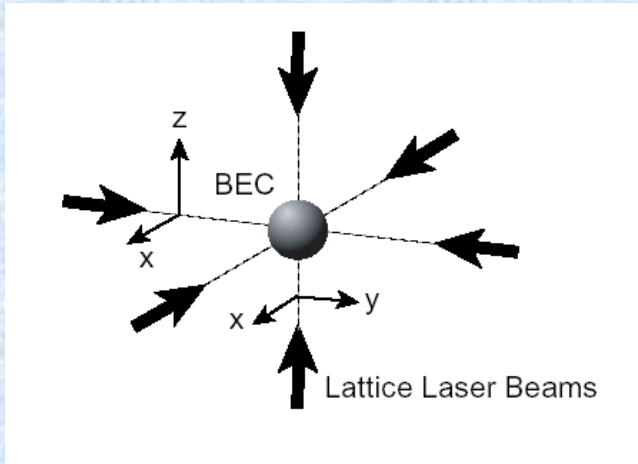


Realization by purpose of a model Hamiltonian (or an effective model) of interest in an experimental setup with highly tunable parameters

Ultracold atoms as quantum simulators of:

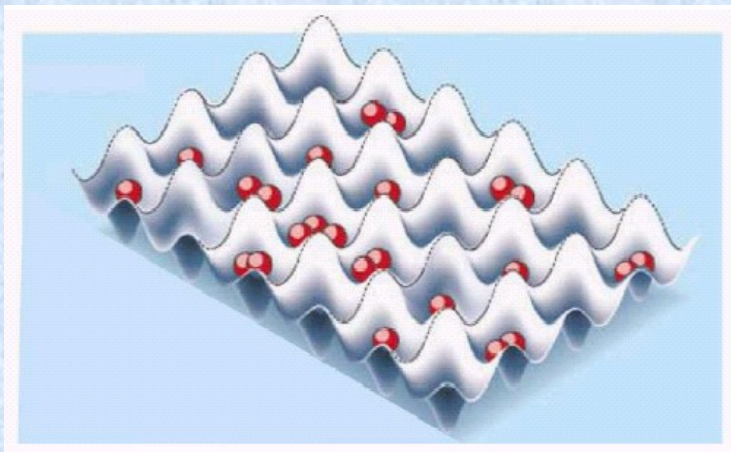
- **Strongly interacting lattice systems (e.g., Fermi and/or Bose Hubbard-like models)**
- **Quantum magnetism**
- **Fermionic superfluids with Cooper pairs**
- **Low-dimensional systems**
- **Quantum Hall physics**
- **Topological states of matter**
- **Field theories**
- **...**

Ultracold bosons in an optical lattice



$$V_{opt}(x) = V_0 \sin^2(kx)$$

e.g., a 1D lattice



It is possible to control:

- barrier height
- interaction term
- the shape of the network
- the dimensionality (1D, 2D, ...)
- the tunneling among planes or among tubes (in order to have a layered structure)

...

Effective Hamiltonian for ultracold bosons in optical lattices (I)

In second quantization, the full quantum many-body Hamiltonian is

$$\hat{H} = \int d\vec{r} \left(\hat{\psi}^\dagger(\vec{r}) \left[\frac{-\hbar^2}{2m} \nabla^2 + V_{opt}(\vec{r}) \right] \hat{\psi}(\vec{r}) + g_0 \hat{\psi}^\dagger(\vec{r}) \hat{\psi}^\dagger(\vec{r}) \hat{\psi}(\vec{r}) \hat{\psi}(\vec{r}) \right)$$

A very good description of (equilibrium and dynamical) low-energy properties - valid for large values of lattice height - is obtained using the Ansatz

$$\hat{\psi}(\vec{r}) = \sum_i \hat{b}_i \Phi_i(\vec{r})$$

tight-binding Ansatz
[D. Jaksch et al., PRL (1998)]

↑
Wannier functions
(to be determined)

One gets...

Effective Hamiltonian for ultracold bosons in optical lattices (II)

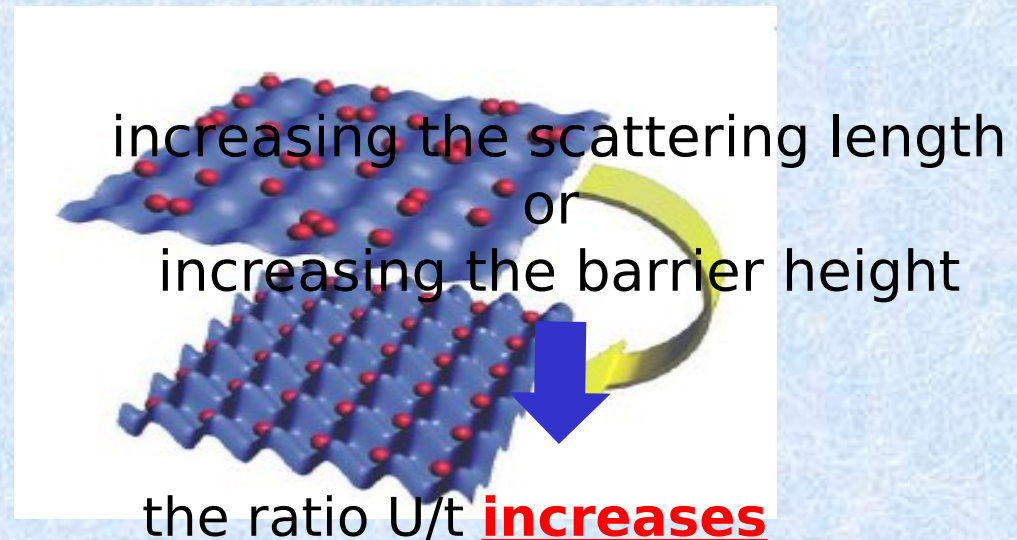
$$\hat{H} = -t \sum_{\langle i, j \rangle} \left(\hat{b}_i^\dagger \hat{b}_j + h.c. \right) + \frac{U}{2} \sum_i \hat{n}_i (\hat{n}_i - 1)$$

$\hat{n}_i \equiv \hat{b}_i^\dagger \hat{b}_i$ N_T number of particles on N sites filling $f = \frac{N_T}{N}$

Bose-Hubbard Hamiltonian

$t/U \gg 1 \rightarrow$ **Superfluid**
dynamics described by the discrete nonlinear Schroedinger equation

$t/U \ll 1 \rightarrow$ **Mott insulator** quantum fluctuations dominate



Effective Hamiltonian for ultracold fermions in optical lattices

Similarly, for a dilute single-species Fermi gas the effective Hamiltonian is

$$\hat{H} = -t \sum_{\langle i, j \rangle} (\hat{c}_i^\dagger \hat{c}_j + h.c.) \equiv -t \sum_{i, j} A_{ij} \hat{c}_i^\dagger \hat{c}_j \quad (f \leq 1)$$

TIGHT-BINDING HAMILTONIAN

Notice that informations about the geometry and the Wannier functions are into the matrix A and the coefficients t, U:

$$t = - \int d\vec{r} \Phi_i^*(\vec{r}) \left[\frac{-\hbar^2}{2m} \nabla^2 + V_{opt}(\vec{r}) \right] \Phi_j(\vec{r})$$
$$U = g_0 \int d\vec{r} |\Phi_i(\vec{r})|^4$$

Quantum simulation of graphene properties (I)



Implementable putting **ultracold fermions** in lattices having **Dirac points**

e.g., in 2D, using the honeycomb lattice itself: using three optical lattices

$$V(x, y) = \sum_{j=1,2,3} V_j \sin^2[k_L(x \cos\theta_j + y \sin\theta_j) + \pi/2]$$

$$\theta_1 = \pi/3, \theta_2 = 2\pi/3, \theta_3 = 0$$



Tight-binding model on the honeycomb
(alias, graphene)

The 3D case

Not a straightforward generalization of the 2D case: indeed, having 2D honeycomb coupled along the z-direction in general destroys the Dirac cones.

More formally:

$$\hat{H} = -t \sum_{i,j} A_{ij} \hat{c}_i^\dagger \hat{c}_j$$

↑
adjacency matrix of the graph [cfr. N. Biggs, Algebraic Graph Theory]

$$-t \sum_j A_{ij} \phi_\alpha(j) = \epsilon_\alpha \phi_\alpha(i) \quad \hat{d}_\alpha = \sum_{\vec{j}} \phi_\alpha(j) \hat{c}_{\vec{j}}$$

$$\hat{H} = \sum_\alpha \epsilon_\alpha \hat{d}_\alpha^\dagger \hat{d}_\alpha$$

The requests are that:

- i) the single particle spectrum has Dirac points (and cones) and the graph has spectral dimension 3**
- ii) that the the adjacency matrix has nearest-neighbour couplings**
- iii) not too many lasers are needed...**

Although symmetries have been studied

[A.A. Abrikosov and S.D. Beneslavskii, JETP (1970) -

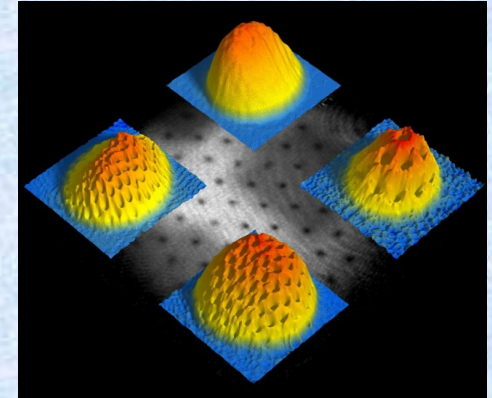
J.L. Manes, PRB (2012)], not easy to satisfy in practice

i)-ii)-iii)...

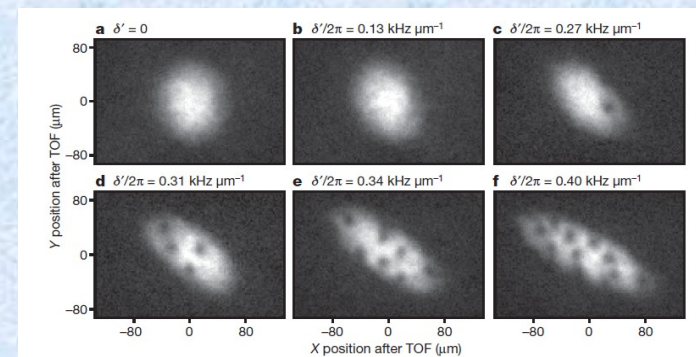
A possible solution: use a synthetic magnetic field

➤ Using rotating traps

[see the review N. Cooper, Adv. Phys. (2008)]



➤ With spatially dependent optical couplings between internal states of the atoms [Y.-J. Lin et al., Nature (2009) - J. Dalibard et al, RMP (2011)]



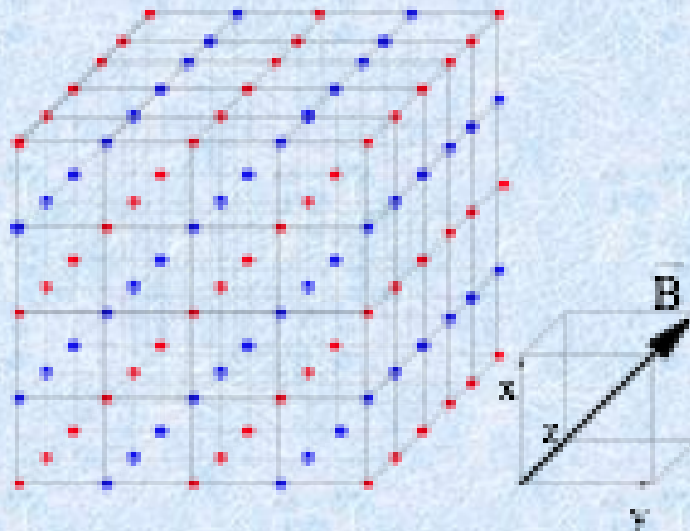
For our purposes: single-species Fermi gas in a π -flux magnetic field (at half filling)

$$\hat{H} = -t \sum_{\langle i, j \rangle} \hat{c}_i^+ e^{-i a_{ij}} \hat{c}_j + h.c.$$

$$a_{ij} = \int_i^j \vec{A} \cdot d\vec{l}$$

$$\vec{B} \equiv \text{rot } \vec{A} = \pi(1, 1, 1)$$

(we can also assume different hoppings t_x , t_y and t_z along the three directions x , y and z)



Single-particle spectrum and Dirac cones (I)

Using the Hasegawa's gauge:

$$\vec{A} = \pi (0, x - y, y - x)$$

[Y. Hasegawa, J. Ph. Soc. Jap. (1990)]

one gets

$$E_{\vec{k}} = \pm 2 \sqrt{t_x^2 \cos^2 k_x + t_y^2 \cos^2 k_y + t_z^2 \cos^2 k_z}$$

with \mathbf{k} belonging to the first (magnetic) Brillouin zone.

[L. Lepori et al, Europhys. Lett. (2010), PRB (2016);
M. Burrello et al, JPA (2017); J. Pinto Barros, M. Burrello, and A. Trombettoni (2020)]

Single-particle spectrum and Dirac cones (II)

- For $t_z = 0$ the results for the 2D case with π -flux are **retrieved** [I. Affleck and J.B. Marston, PRB (1988)] **are retrieved**.
- Excitations around the two inequivalent Dirac points **obey the 3D Dirac equation**.
- In the limit of vanishing t_z one retrieves the 2D Dirac equation.
- A mass term can be added using a Bragg pulse.
- The Dirac points does not depend on t_x , t_y and t_z .
- With an attractive interaction U one has a semimetal-superfluid transition at a finite value of U
- Extendable to many components

Applications

- With a spatial control of the synthetic magnetic field one can an e.m. field.
- With a dynamical gauge field one can then have a simulation of the 3+1 QED.
- One may also think to have the fermions living in one dimension and the gauge field in an higher dimension: this has been studied in the context of graphene, giving rise to pseudo-QED [E.C. Marino, Nucl. Phys. B, 1993]. In one dimension one would then have the pseudo-Schwinger model.

Outline

- Ultracold atoms as quantum simulators for relativistic field theories
- **Symmetry-locked superfluid phases**
- Topological Kondo model in junctions of Tonks-Girardeau gases

Small Dictionary (I)

Abelian transformation: only
one parameter/generator

$$U = e^{i\phi}$$

Non Abelian symmetry : more
than one parameter/generator,
not commuting each others

$$U = e^{i\alpha_j T^j}$$

$$[T_i, T_k] \neq 0$$

Small Dictionary (II)

Global transformation :
parameters **NOT** space-time
dependent

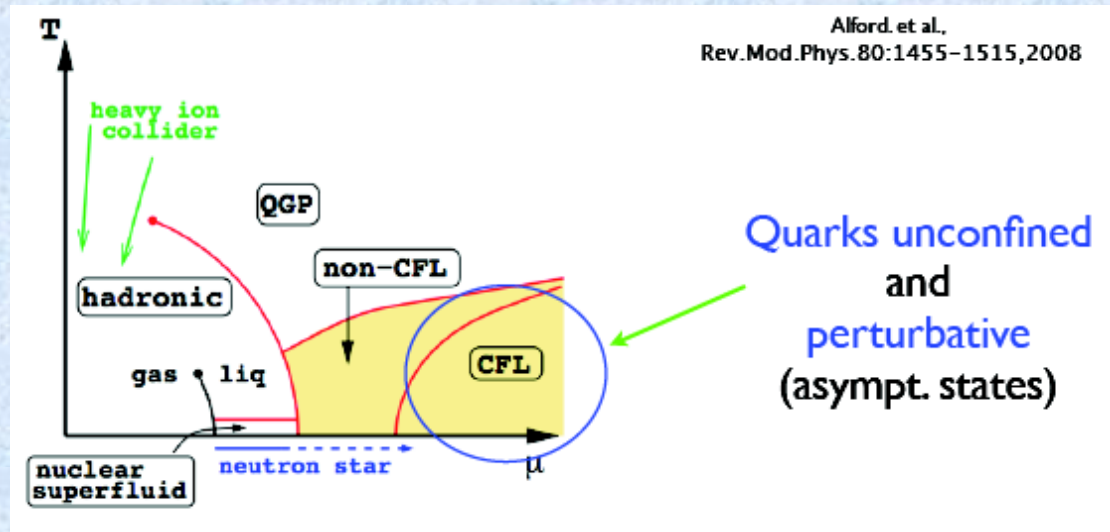
$$U = e^{ia_j T^j}$$

Local (gauge) transformation :
parameters space-time
dependent

$$U(x^\mu) = e^{ia_j(x^\mu) T^j}$$

Motivations

- physics of interacting fermionic mixtures
- QCD inspired problem(s)



The proposed quantum simulation

→ 4 fermionic species: e.g., 4 species of Yb or a mixture ^{171}Yb - ^{173}Yb

→ 2 doublets $\begin{pmatrix} c_1 \\ c_2 \end{pmatrix}$ $\begin{pmatrix} f_1 \\ f_2 \end{pmatrix}$

→ $U(2)_c \times U(2)_f$ (global) invariant Hamiltonian


→
$$\hat{H}_{int} = -U_c \sum_{i,c} (c_{i;c}^\dagger c_{i;c})^2 - U_f \sum_{i,f} (c_{i;f}^\dagger c_{i;f})^2 +$$
$$-U_{cf} \sum_{i,c,f} c_{i;c}^\dagger c_{i;c} c_{i;f}^\dagger c_{i;f}$$

$$\begin{pmatrix} c_1 \\ c_2 \end{pmatrix} \equiv \begin{pmatrix} c_g \\ c_r \end{pmatrix}$$

$$\begin{pmatrix} f_1 \\ f_2 \end{pmatrix} \equiv \begin{pmatrix} c_u \\ c_d \end{pmatrix}$$

Order parameters

$$\langle c_i; u c_i; d \rangle \equiv \Delta_f, \quad \langle c_i; r c_i; g \rangle \equiv \Delta_c, \quad \langle c_i; c c_i; f \rangle \equiv \Delta_{cf}$$


a 2x2 matrix

Order parameters:

$$2|\Delta_0|^2 = (|\Delta_c|^2 + |\Delta_f|^2)$$

$$\Delta_+^2 = \text{Tr} \left(\Delta_{cf}^\dagger \Delta_{cf} \right), \quad \Delta_-^2 = 2 \det \Delta_{cf}$$

We consider $U_c = U_f \equiv U$

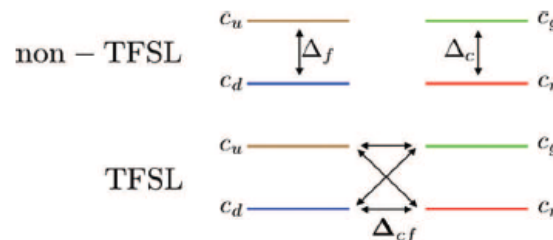
Results (I)

$$2|\Delta_0|^2 = (|\Delta_c|^2 + |\Delta_f|^2)$$

$$\Delta_+^2 = \text{Tr} \left(\Delta_{cf}^\dagger \Delta_{cf} \right), \quad \Delta_-^2 = 2 \det \Delta_{cf}$$

We find a “two-flavors” symmetry-locked phase (TFSL) for $U_{cf} > U$:

The minimization of F with respect to Δ_\pm and Δ_0 gives $|\Delta_+| = |\Delta_-|$ and $|\Delta_c| = |\Delta_f|$. We find that for $U_{cf} \neq U$ the gap equations are not consistent if both Δ_+ and Δ_0 are non-zero both $T = 0$ and finite temperature and two phases are found as follows (see fig. 1): i) Non-TFSL phase: for $U_{cf} < U$ it is $\Delta_+ = 0$ and $\Delta_0 \neq 0$; ii) TFSL phase: for $U_{cf} > U$ it is $\Delta_0 = 0$ and $\Delta_+ \neq 0$.



Results (II)

Non-TFSL superfluid phase **abelian**:

$$U(2)_c \times U(2)_f \rightarrow SU(2)_c \times SU(2)_f$$

“Two-Flavours locking” phase :

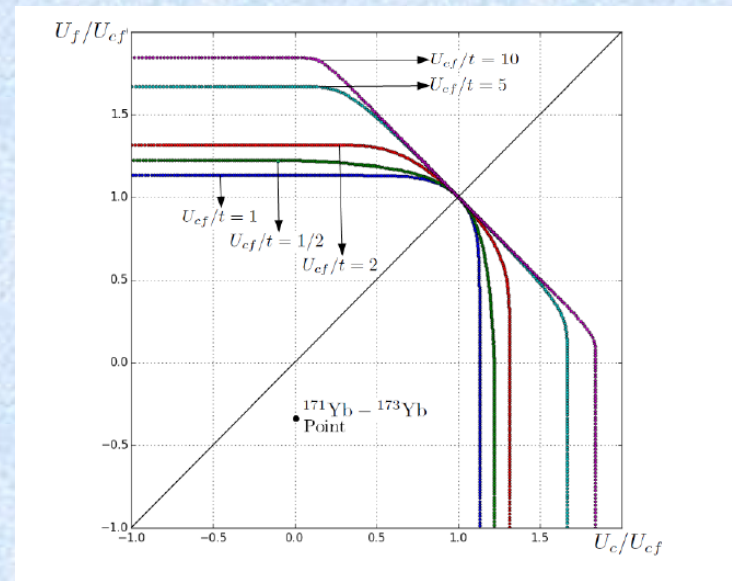
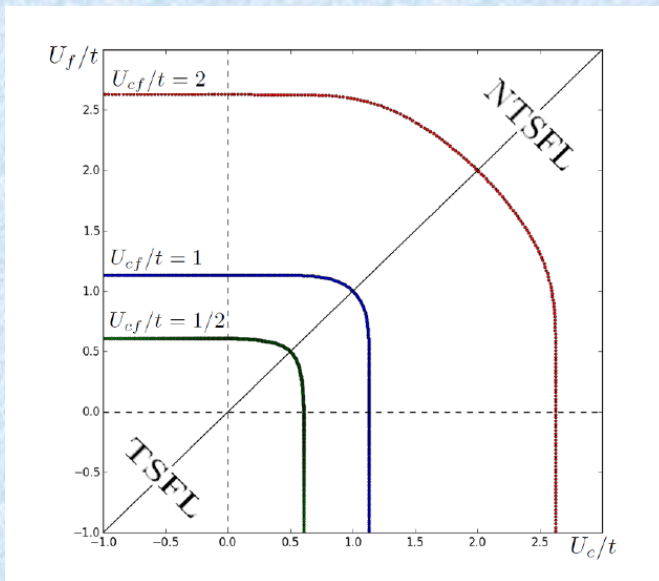
Spont. Symm. Break. $U(2)_c \times U(2)_f \rightarrow U(2)_{c+f}$



vortices in the CFL phase have fractional flux

Results (III)

But it does survive when interactions are different? Especially if $U_f < 0$...



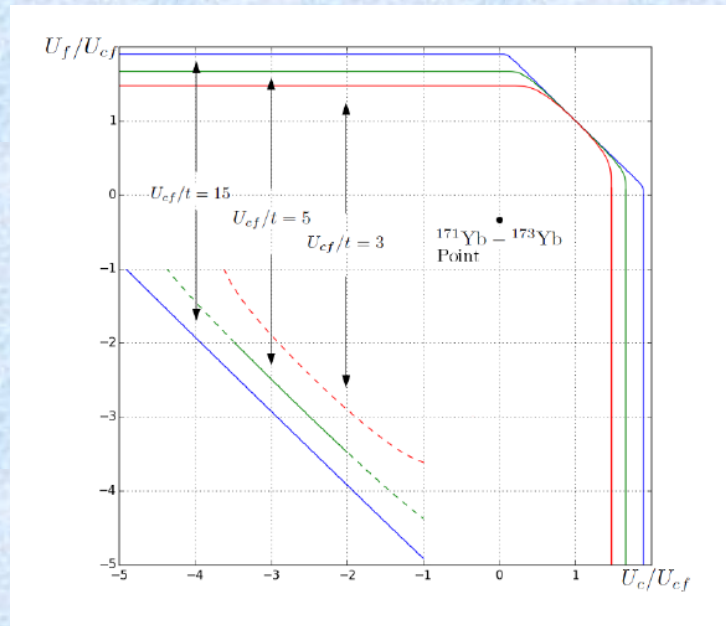
$$a_{171-173} = -578a_0$$

$$a_{173-173} = +200a_0$$

$$a_{171-171} = -3a_0$$

Results (IV)

The previous are mean-field results → combining with a strong-coupling computation



To have a “true” color-flavor locking: even without putting interactions, use dynamical gauge fields

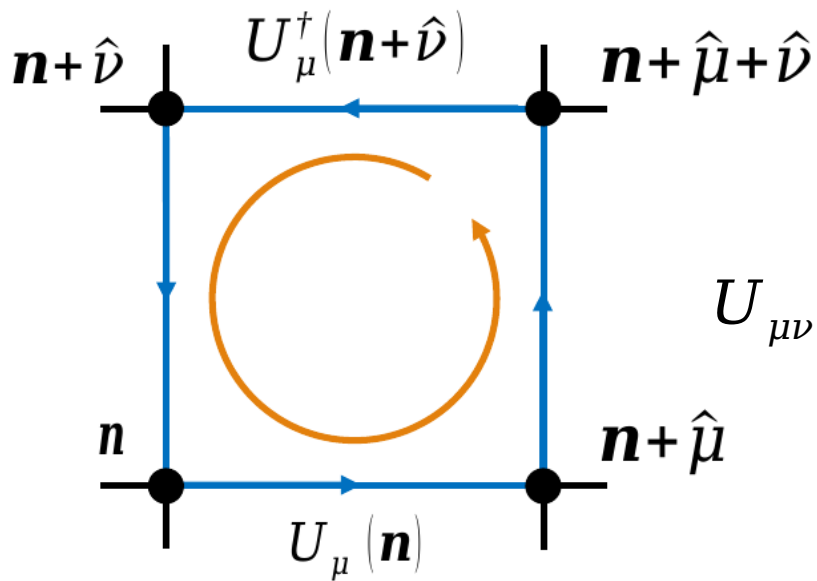
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- Symmetry-locked superfluid phases
- **Link models → how to generate plaquette terms?**

Basic quantities on the lattice



[Wilson, PRD (1974)]

$$U_{\mu}(\mathbf{n}) = e^{ieaA_{\mu}(\mathbf{n})}$$

LINKS

$$U_{\mu\nu}(\mathbf{n}) = U_{\mu}(\mathbf{n})U_{\nu}(\mathbf{n} + \hat{\mu})U_{\mu}^+(\mathbf{n} + \hat{\nu})U_{\nu}^+(\mathbf{n})$$

$$U_{\mu\nu}(\mathbf{n}) = e^{iea^2F_{\mu\nu}(\mathbf{n})}$$

PLAQUETTES

Link models

ELECTRIC FIELD

$$[U_\mu(\mathbf{n}), E_\nu(\mathbf{n}')] = -\delta_{\mu,\nu}\delta_{\mathbf{n},\mathbf{n}'}U_\mu(\mathbf{n})$$

KOGUT-SUSSKIND HAMILTONIAN

$$H_g = \frac{e^2}{2} \sum_{\mathbf{n},\mu} E_\mu^2(\mathbf{n}) - \frac{1}{4a^2 e^2} \sum_P (U_{\mu\nu} + U_{\mu\nu}^\dagger)$$

ROKHSAR-KIVELSON HAMILTONIAN

$$H_{RK} = H_g + \lambda \sum_P (U_{\mu\nu} + U_{\mu\nu}^\dagger)^2$$

Quantum link models → replace the Wilson operators by discrete quantum degrees of freedom, still living on the links of the lattice (quantum links)

Link models (II)

Bosonic quantum link models

$$U_{\mu}(\mathbf{n}) = S_{\mu}^{+}(\mathbf{n}), \quad U_{\mu}^{\dagger}(\mathbf{n}) = S_{\mu}^{-}(\mathbf{n}), \quad E_{\mu}(\mathbf{n}) = S_{\mu}^{z}(\mathbf{n})$$

Fermionic quantum link models: in terms of fermionic states and occupation numbers, we denote the two states of the local Hilbert space with $|0\rangle$ and $|1\rangle$

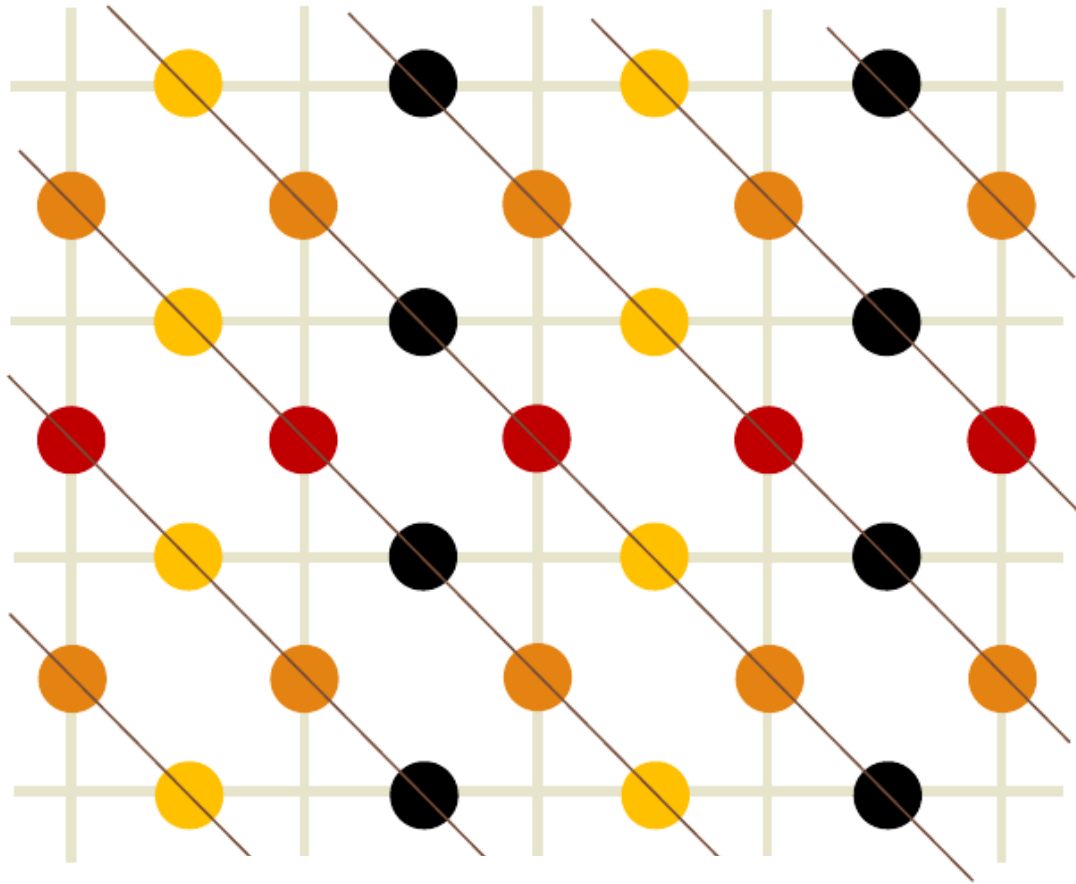
$$|1\rangle = c_{\mu}^{\dagger}(\mathbf{n})|0\rangle \quad U_{\mu\nu}(\mathbf{n}) = c_{\mu}(\mathbf{n})c_{\nu}(\mathbf{n} + \hat{\nu})c_{\mu}^{\dagger}(\mathbf{n} + \hat{\nu})c_{\nu}^{\dagger}(\mathbf{n})$$

Proposals for having plaquette terms

→ 4 correlated hoppings + angular momentum conservation
[Zohar, Cirac, and Reznik, PRA (2013)]

→ Dual formulation: single hopping + conditional operations on the
nearest-neighbours [A. Celi et al, PRX (2020)]

A proposal using a spin dependent-optical lattice



$$H_0 = -h \sum_{x,m} \epsilon_{xm} b_{xm}^\dagger b_{xm}$$

Derivation of the plaquette term in perturbation theory

$$H_1 = H_{\text{hop}} + H_{\text{int}} \equiv -t \sum_{\langle i,j \rangle_d, m} (b_{im}^\dagger b_{jm} + \text{h.c.}) + \frac{1}{2} \sum_{\langle i,j \rangle, m, m'} V_{mm'}^{i,j} b_{xm}^\dagger b_{ym'}^\dagger b_{xm'} b_{ym}$$

(hard-core condition assumed)

At the third order of the perturbation theory for large \hbar one finds:

$$H^{(\text{eff})} = \frac{t^2}{\hbar} \sum_{\langle i,j \rangle_d, m, m'} n_{im} n_{jm'} - \frac{1}{\hbar} \sum_{\langle i,j \rangle, m, m'} (V_{mm'}^{i,j})^2 n_{im} n_{jm'} + \frac{t^2}{\hbar} \sum_{\substack{i, i', j, j' \in \square \\ m, m'}} V_{mm'}^{(i, i')} b_{j'm}^\dagger b_{jm'}^\dagger b_{i'm'} b_{im}$$

[P. Fontana, J. Pinto-Barros, M. Burrello, and A. Trombettoni, to be submitted]

Connections with link models

$$U_{im} = b_{im}^\dagger, \quad U_{im}^\dagger = b_{im}$$

$$E_{im} \equiv n_{im} - \frac{1}{2}$$

$$H^{(\text{eff})} = \sum_{\langle x,y \rangle_d, m, m'} \lambda_1^{(mm')} E_{xm} E_{ym'} - \sum_{\langle x,y \rangle, m, m'} \lambda_2^{(mm')} E_{xm} E_{ym} - J \sum_{\square} (U_{\square} + U_{\square}^\dagger)$$

Conclusions

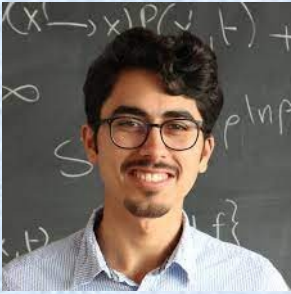
Ultracold atoms as quantum simulators for field theories models

Ultracold fermions and gauge fields are a tool to emulate mechanisms such as the color-flavor locking

Quantum simulation of link models with plaquette terms is demanding → we discussed a proposal involves spinor dipolar gases

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