



Overview of QIS and Quantum Emulation studies at University of Illinois

>> *Prospects for studying dynamical gauge fields with cold atoms and molecules*

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University of Illinois at Urbana-Champaign

Intersections between Nuclear Physics and Quantum Information (NPQI workshop)
Argonne National Laboratory – 2 pm, March 29, 2018

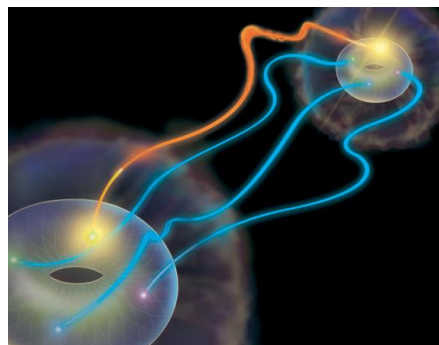
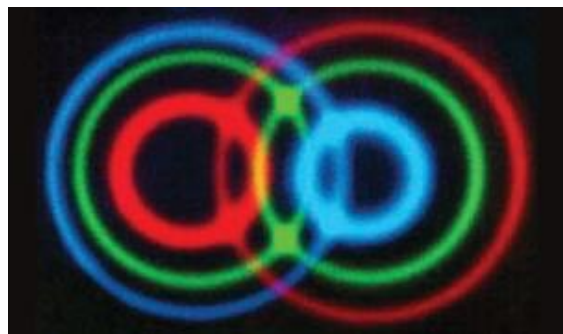
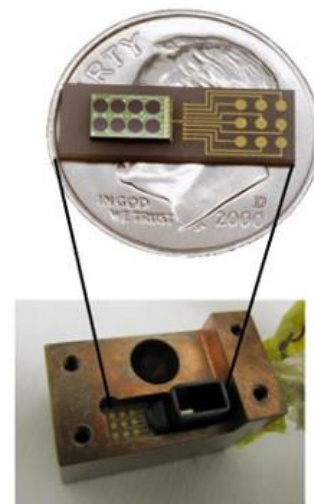
Kwiat Group: Quantum Information Science

Improving quantum technologies

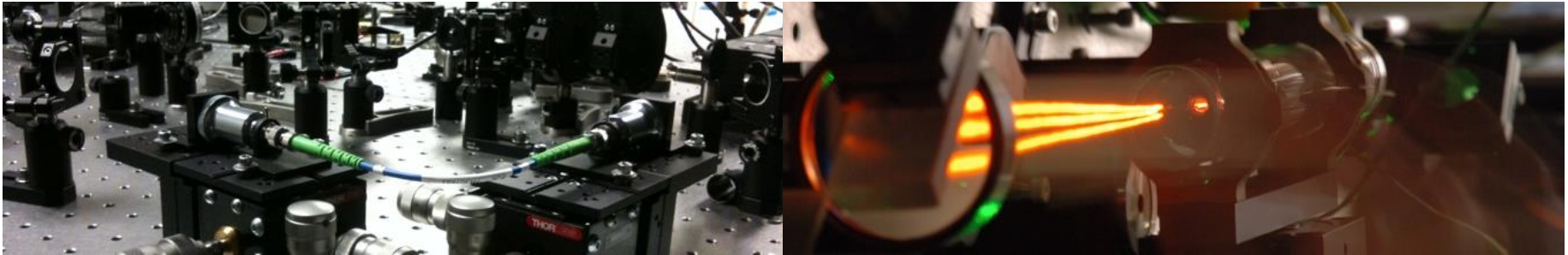
- Development of entangled photon and single-photon sources
- Development of single-photon detectors
- Improving efficiency of photon collection in noisy environments

More fundamental studies

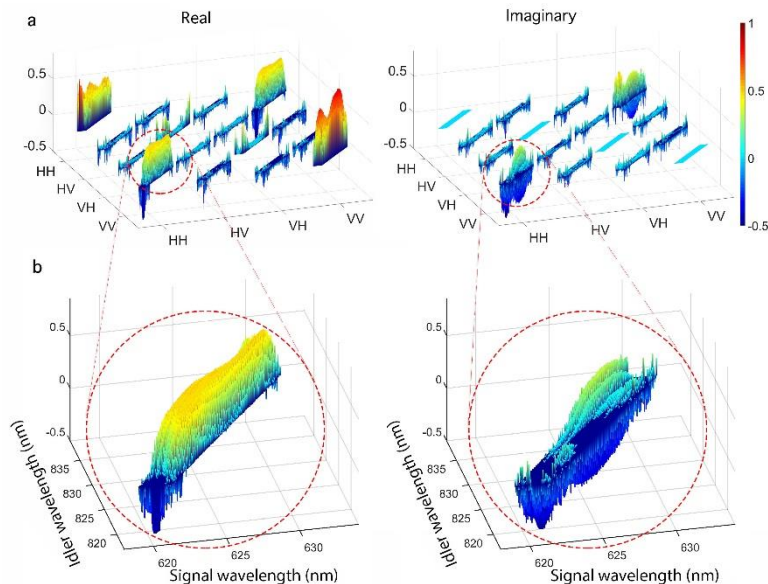
- Bell-like tests of nonlocal realism
- Superdense encoding of quantum information
- Teleportation and quantum communications
- Few- and single-photon detection by humans



Lorenz Group: Quantum Optics and Spectroscopy



Motivation: quantum information applications, fundamental studies

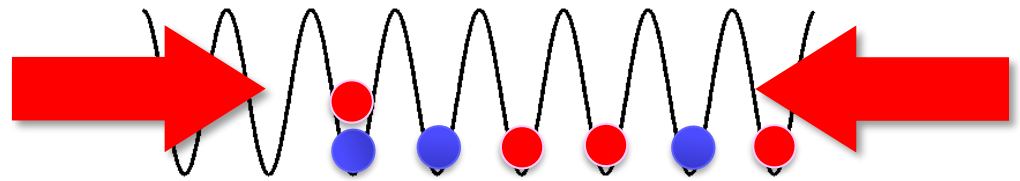
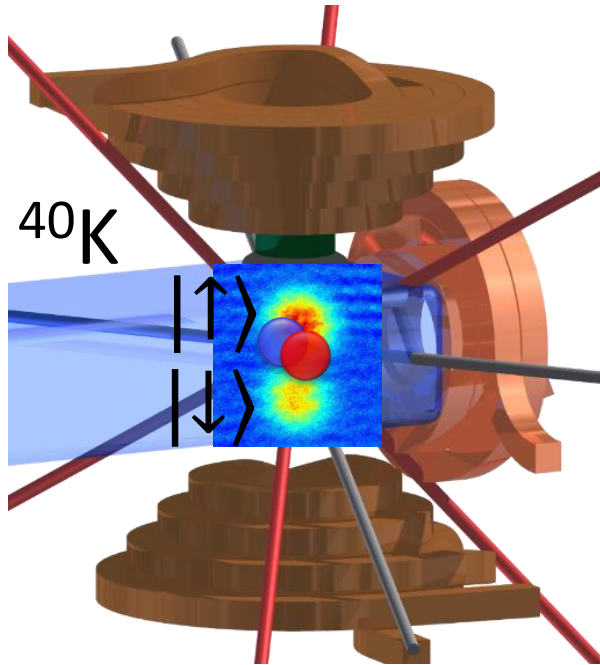


- Engineering photonic quantum states for quantum applications
- High-resolution, high-efficiency measurement techniques for characterization of photonic quantum states in multiple degrees of freedom
- Quantum memories for telecom broadband photonic quantum states, fundamental studies of influence of light-matter interaction on photon-pair properties



DeMarco group

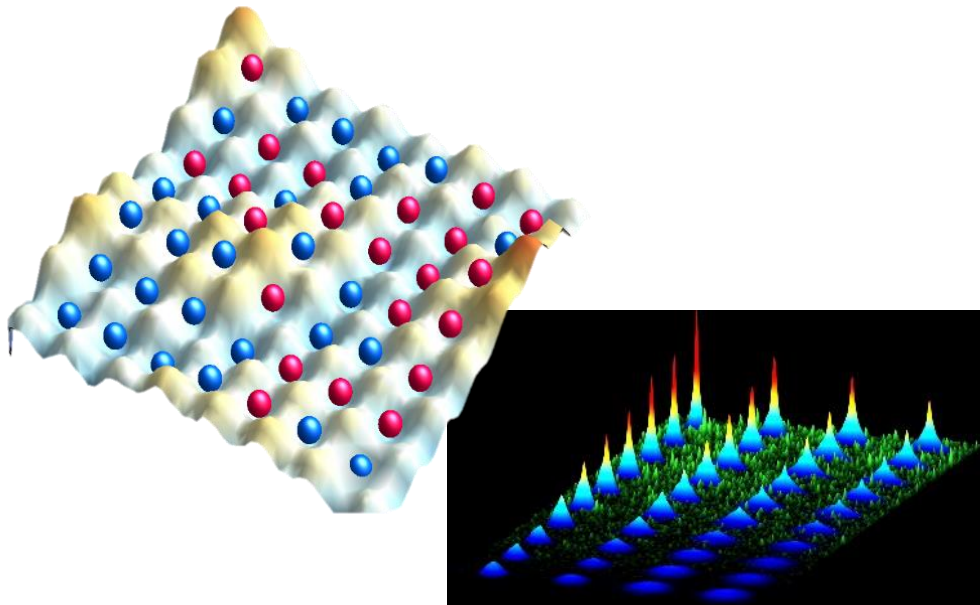
Optical lattice quantum emulation



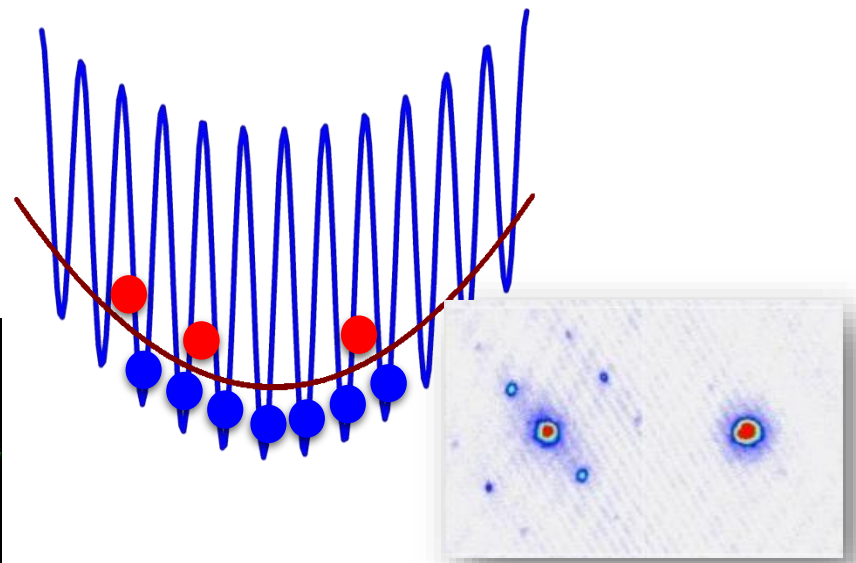
$$H = -t \sum_{\langle ij \rangle} c_i^\dagger c_j + U \sum_i n_{\downarrow i} n_{\uparrow j}$$

Hubbard model

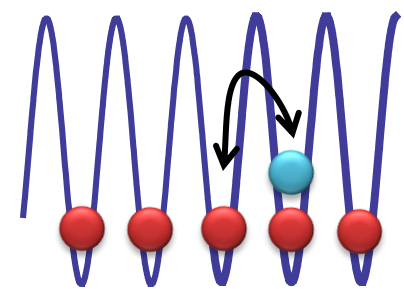
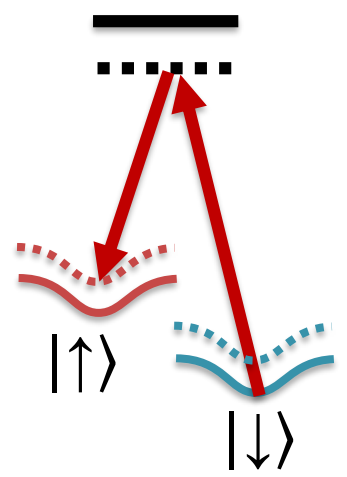
Minimal model for strongly interacting quantum matter



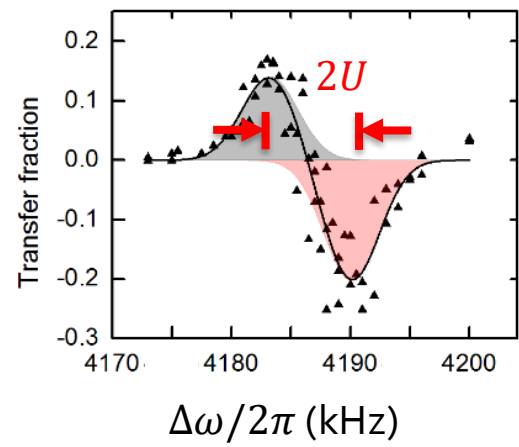
Disorder



Spin-dependent lattices & potentials



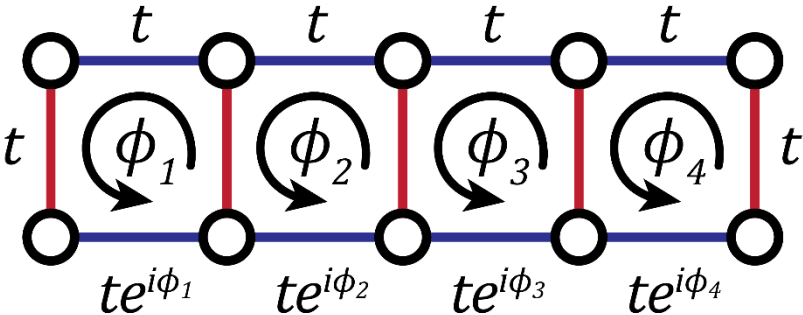
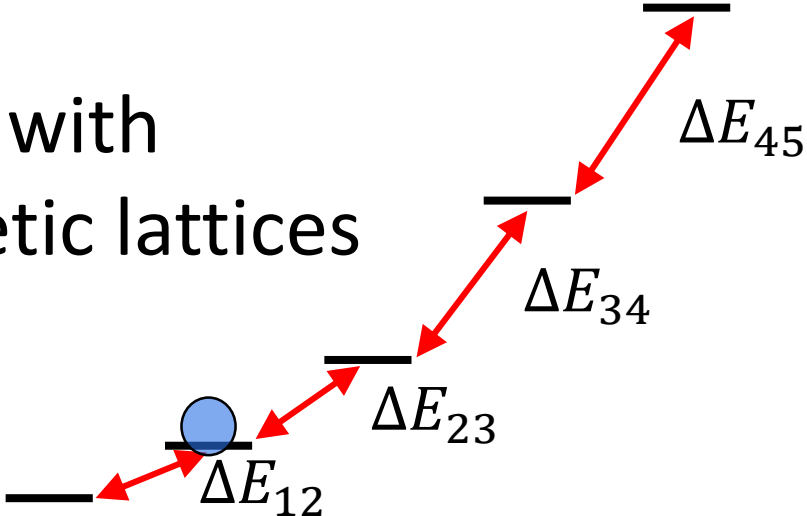
Correlated Tunneling



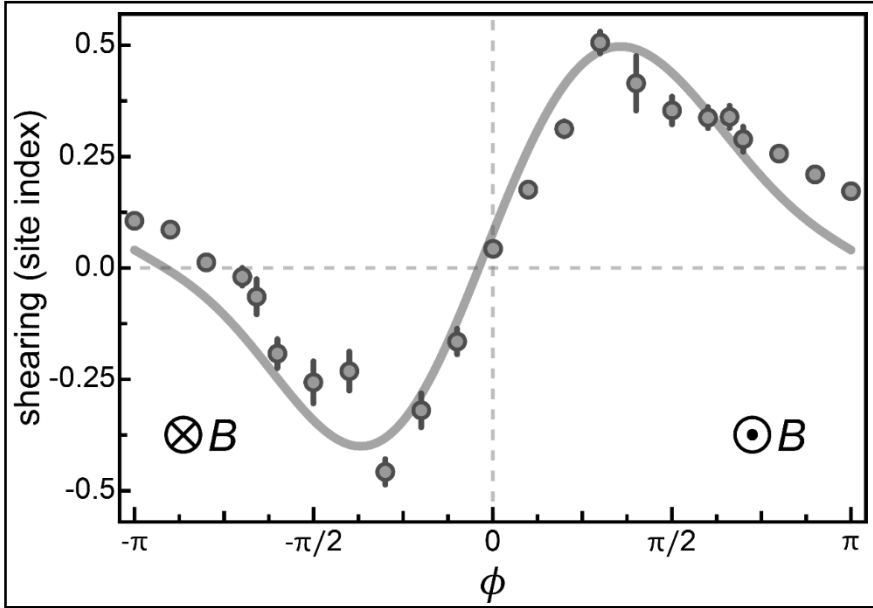
Gadway group

Quantum simulation with
non-standard, synthetic lattices

w/ cold atoms & molecules

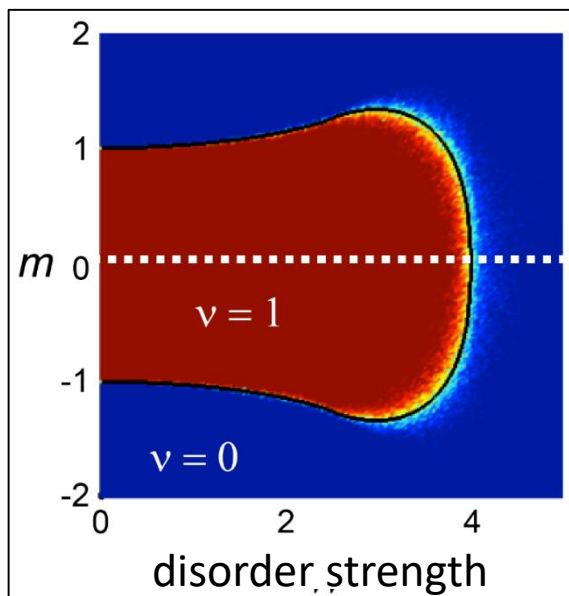


Designer *classical* gauge fields



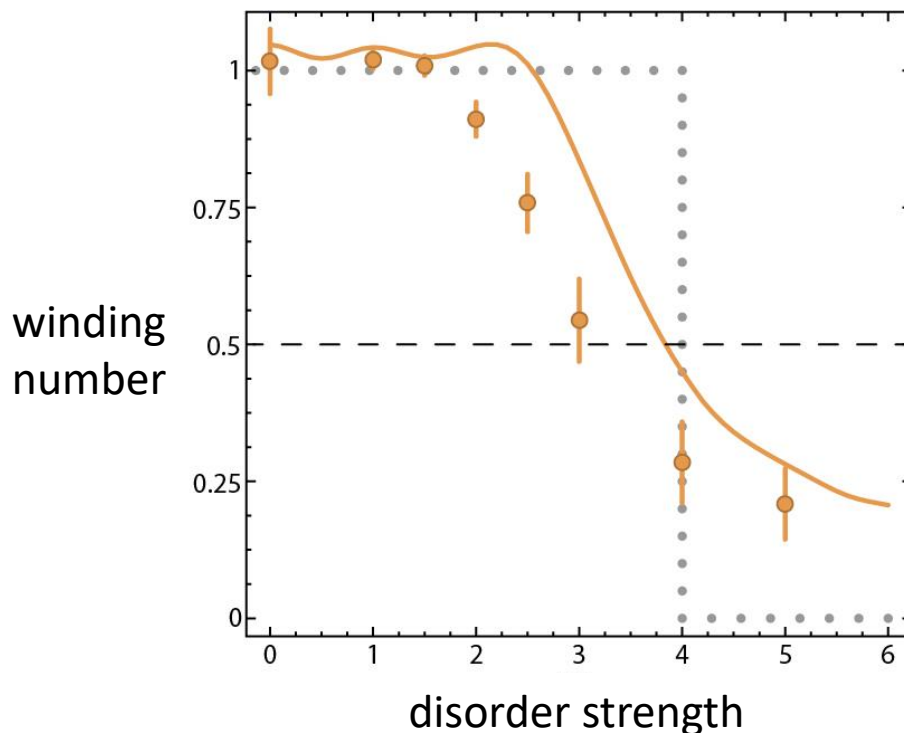
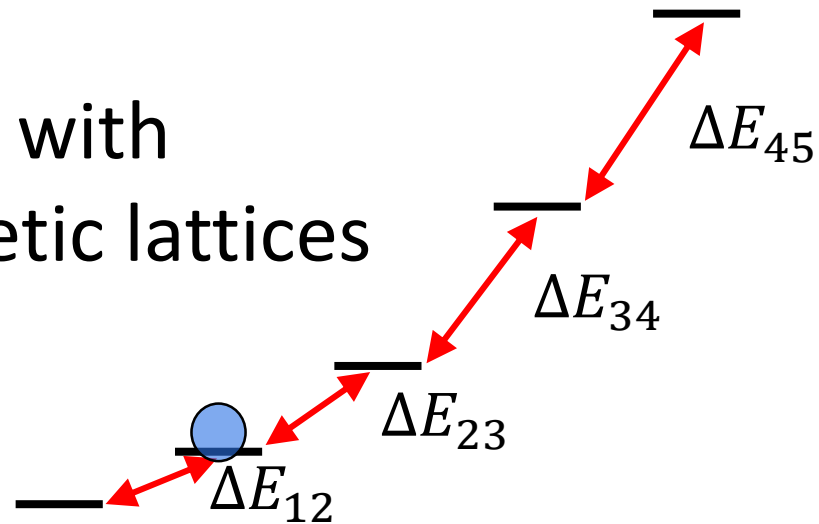
Gadway group

Quantum simulation with
non-standard, synthetic lattices



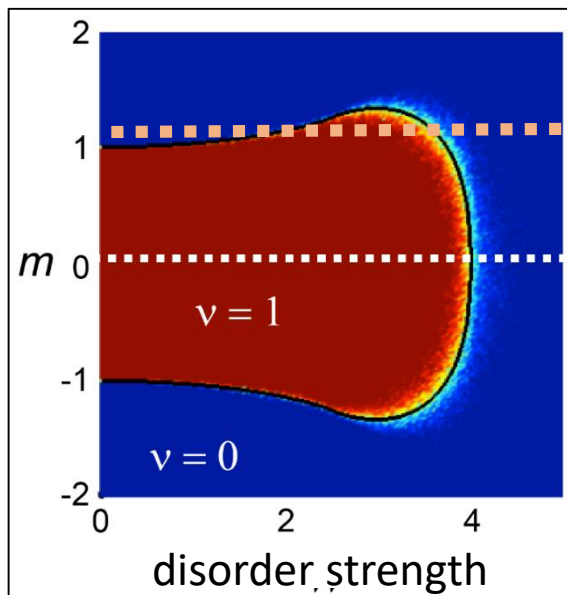
Interplay of:

- topology & disorder
- topology & interactions
- disorder & interactions



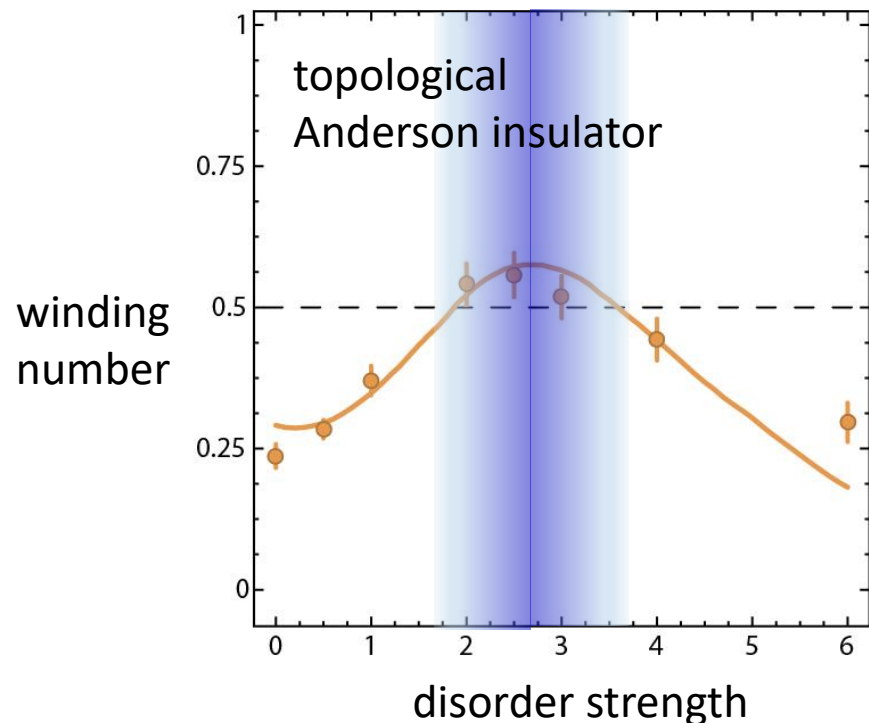
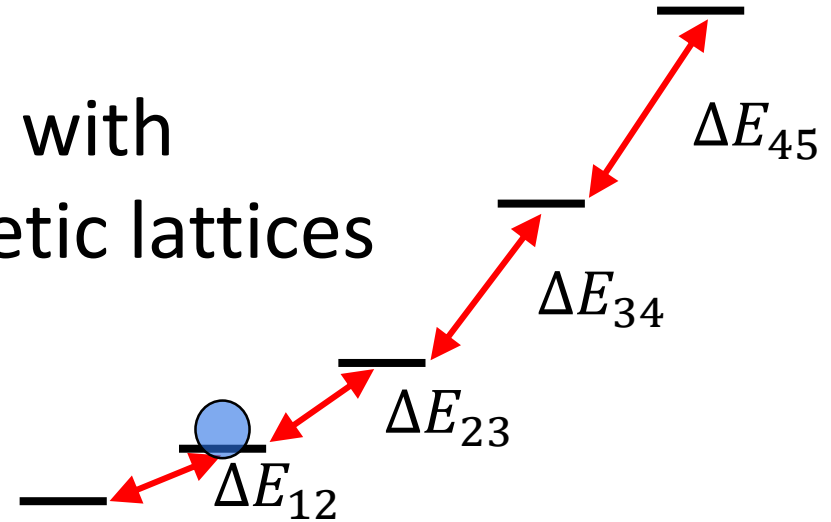
Gadway group

Quantum simulation with
non-standard, synthetic lattices

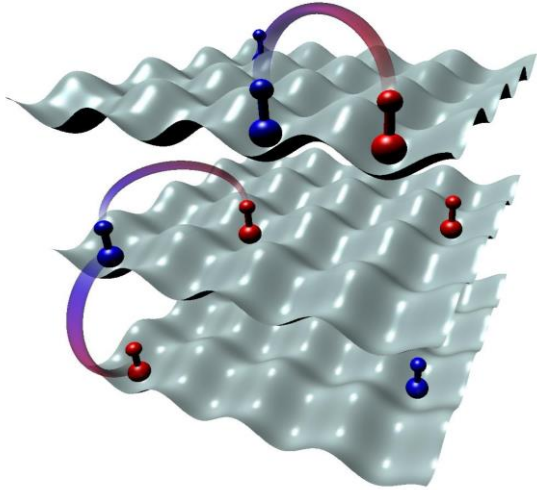


Interplay of:

- topology & disorder
- topology & interactions
- disorder & interactions



Gadway / DeMarco – ultracold molecules



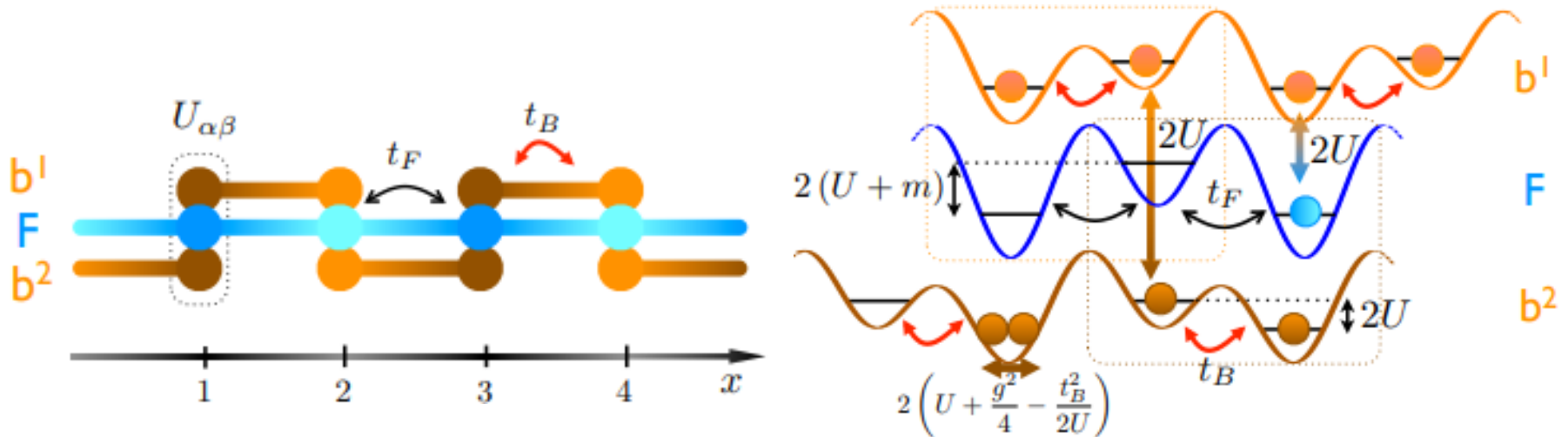
Harnessing dipolar exchange interactions
for the emulation of lattice gauge theories

Gadway / DeMarco – ultracold molecules

Requires extreme control, not very forgiving for cold atoms
→ molecules are a natural fit

Quantum field theory
Lattice gauge theory
Quantum link model

- Multiple species (2 bosons, 1 fermion)
- Spin-dependent tunneling / site energies
- Control of all interactions [must be stable]
- Keep entropy low
→ gauge invariance only at low T, low S/N

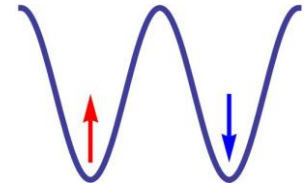


Banerjee, et al. PRL 109, 175302 (2012)

Emulating LGTs with ultracold molecules

atoms \rightarrow molecules

particle hopping \rightarrow spin “hopping” by dipolar exchange



correlated hopping (low S/N , low t/U) \rightarrow naturally imposed by dipolar exchange

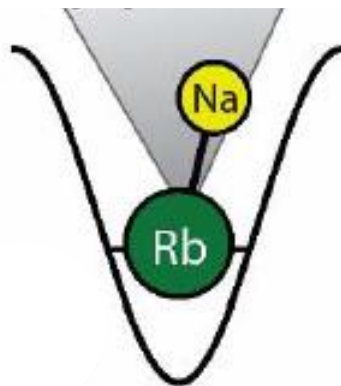
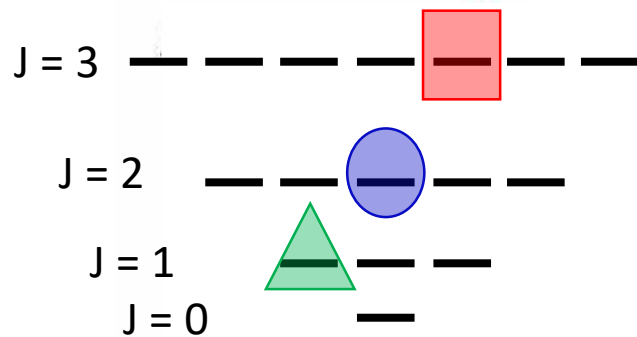
spin-dependent potentials \rightarrow local “magnetic fields” / light shifts

start with “low” entropy by motion \rightarrow motion frozen out, initial spin entropy = 0

gauge invariance / Gauss’ law imposed by hand

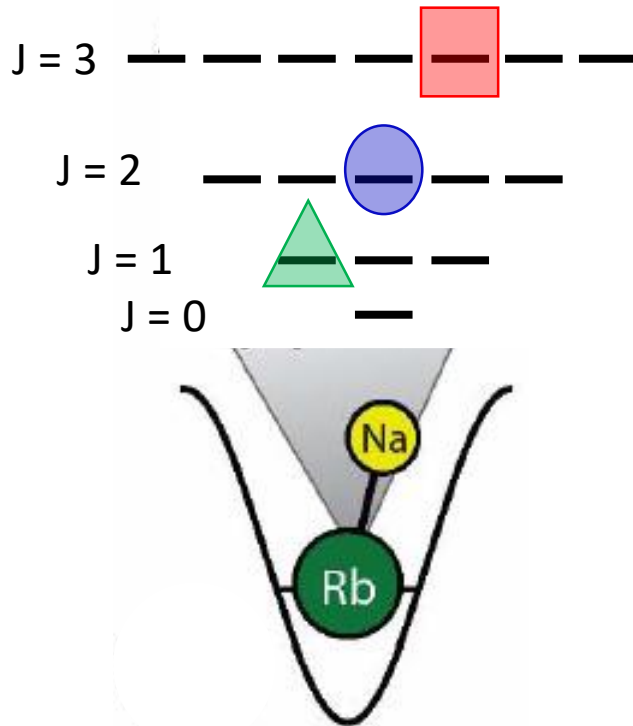
Emulating LGTs with ultracold molecules

internal rotational states

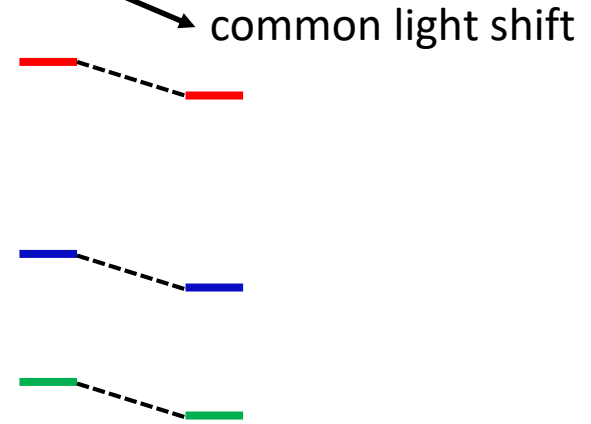


Emulating LGTs with ultracold molecules

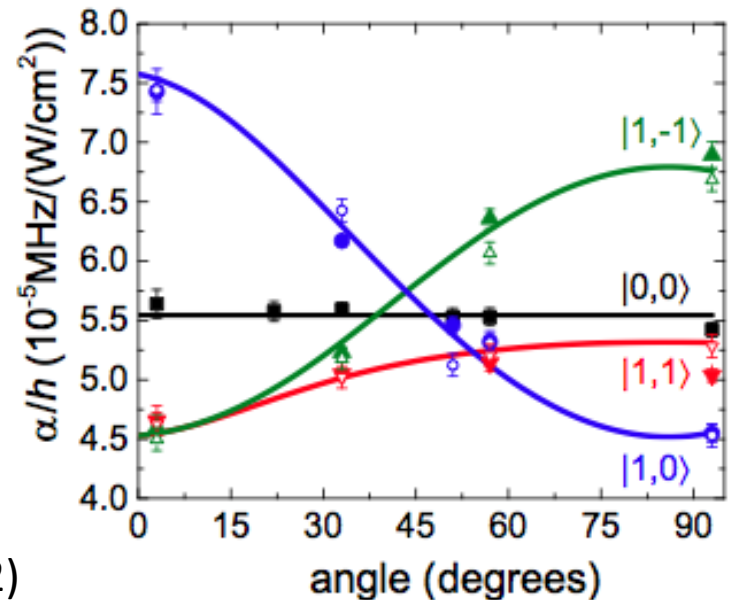
internal rotational states



free space

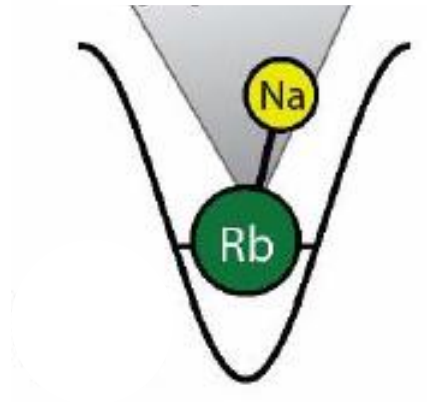
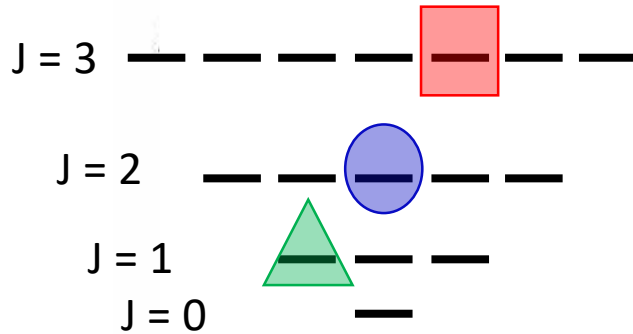


anisotropic polarizability
Neyenhuis, *et al.* PRL (2012)

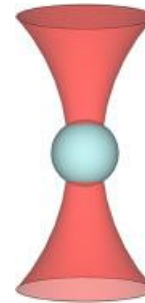
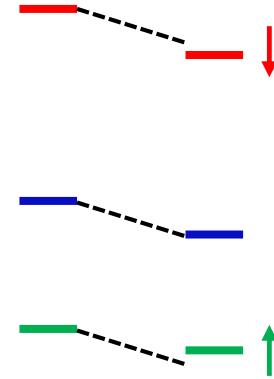


Emulating LGTs with ultracold molecules

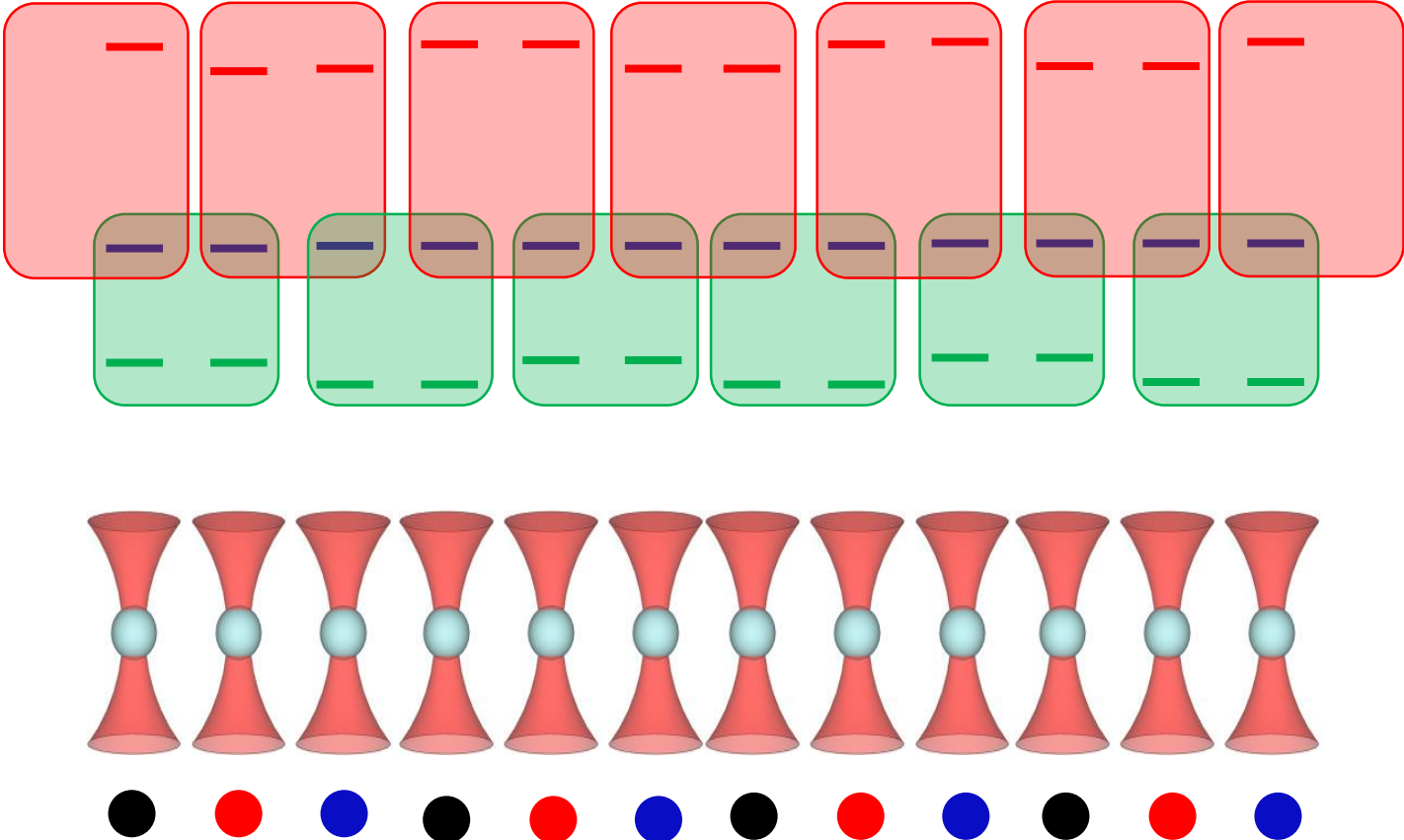
internal rotational states



free space (I, \hat{e})
state-dep. light shift

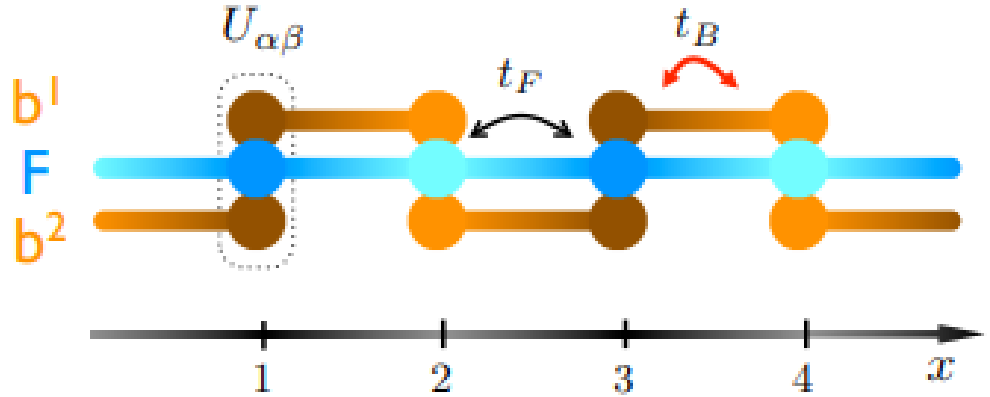
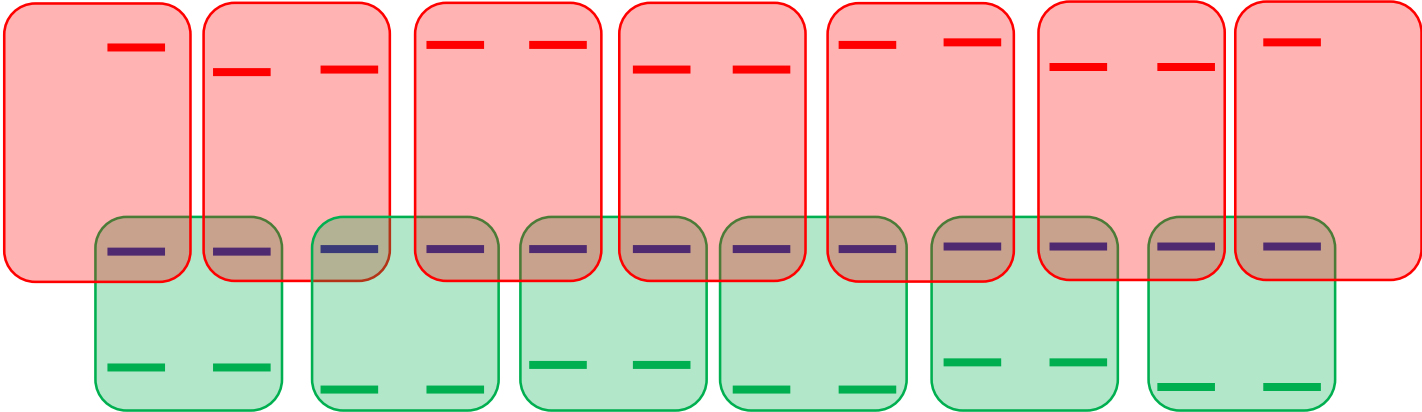


Emulating LGTs with ultracold molecules



$$(I_1, \hat{e}_1) , (I_2, \hat{e}_2) , (I_3, \hat{e}_3)$$

Emulating LGTs with ultracold molecules



Banerjee, et al. PRL 109, 175302 (2012)

Thanks! Questions?

