

Quantum Information Processing and Quantum Simulation with Ultracold Alkaline-Earth Atoms in Optical Lattices

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Quantum Simulation: *Nature Phys.* 6, 289 (2010)



KITP

November 19, 2010

Alkaline-earth(-like) atoms

| | | | | | | | | | | | | | | | | | | | | | | | |
|--------------------------------|---------------------------------|---------------------------------|---------------------------------|-------------------------------------|---------------------------------|----------------------------------|---------------------------------|-------------------------------|----------------------------------|------------------------------------|--------------------------------|---------------------------------|------------------------------------|--------------------------------|---------------------------------|-------------------------------|------------------------------|-----------------------------|----------------------------|------------------------------|----------------------------|------------------------------|----------------------------|
| hydrogen 1 H 1.0079 | beryllium 4 Be 9.0122 | | | | | | | | | | | | | | | | | helium 2 He 4.0026 | | | | | |
| lithium 3 Li 6.941 | magnesium 12 Mg 24.305 | | | | | | | | | | | | | | | | | boron 5 B 10.811 | carbon 6 C 12.011 | nitrogen 7 N 14.007 | oxygen 8 O 15.999 | fluorine 9 F 18.998 | neon 10 Ne 20.180 |
| sodium 11 Na 22.990 | calcium 20 Ca 40.078 | scandium 21 Sc 44.956 | titanium 22 Ti 47.867 | vanadium 23 V 50.942 | chromium 24 Cr 51.996 | manganese 25 Mn 54.938 | iron 26 Fe 55.845 | cobalt 27 Co 58.933 | nickel 28 Ni 58.693 | copper 29 Cu 63.546 | zinc 30 Zn 65.39 | gallium 31 Ga 69.723 | germanium 32 Ge 72.61 | arsenic 33 As 74.922 | selenium 34 Se 78.96 | bromine 35 Br 79.904 | krypton 36 Kr 83.80 | | | | | | |
| potassium 19 K 39.098 | strontium 38 Sr 87.62 | yttrium 39 Y 88.906 | zirconium 40 Zr 91.224 | niobium 41 Nb 92.906 | molybdenum 42 Mo 95.94 | technetium 43 Tc [98] | ruthenium 44 Ru 101.07 | rhodium 45 Rh 102.91 | palladium 46 Pd 106.42 | silver 47 Ag 107.87 | cadmium 48 Cd 112.41 | indium 49 In 114.82 | tin 50 Sn 118.71 | antimony 51 Sb 121.76 | tellurium 52 Te 127.60 | iodine 53 I 126.90 | xenon 54 Xe 131.29 | | | | | | |
| rubidium 37 Rb 85.468 | barium 56 Ba 137.33 | lanthanum 57 La 138.91 | hafnium 72 Hf 178.49 | tantalum 73 Ta 180.95 | tungsten 74 W 183.84 | rhenium 75 Re 186.21 | osmium 76 Os 190.23 | iridium 77 Ir 192.22 | platinum 78 Pt 195.08 | gold 79 Au 196.97 | mercury 80 Hg 200.59 | thallium 81 Tl 204.38 | lead 82 Pb 207.2 | bismuth 83 Bi 208.98 | polonium 84 Po [209] | astatine 85 At [210] | radon 86 Rn [222] | | | | | | |
| caesium 55 Cs 132.91 | radium 88 Ra [226] | actinium 87 Ac [227] | lutetium 71 Lu 174.97 | rutherfordium 104 Rf [261] | dubnium 105 Db [262] | seaborgium 106 Sg [266] | bohrium 107 Bh [264] | hassium 108 Hs [269] | meitnerium 109 Mt [268] | ununnillium 110 Uun [271] | ununium 111 Uuu [272] | ununbium 112 Uub [277] | ununquadium 114 Uuq [289] | | | | | | | | | | |

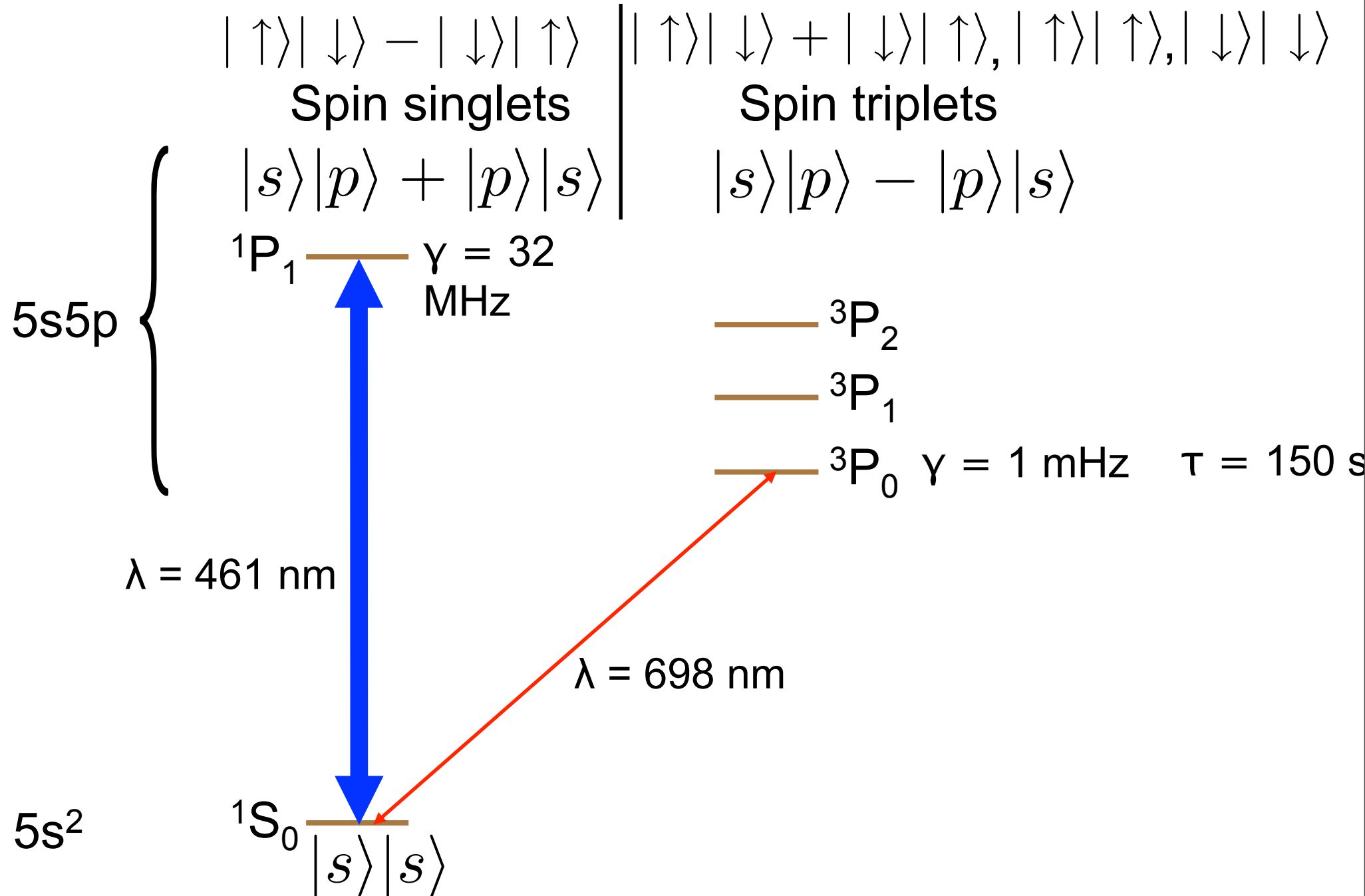
* Lanthanide series

** Actinide series

| | | | | | | | | | | | | | |
|---------------------------------|-------------------------------|------------------------------------|---------------------------------|---------------------------------|--------------------------------|--------------------------------|----------------------------------|--------------------------------|----------------------------------|----------------------------------|-------------------------------|-----------------------------------|---------------------------------|
| lanthanum 57 La 138.91 | cerium 58 Ce 140.12 | praseodymium 59 Pr 140.91 | neodymium 60 Nd 144.24 | promethium 61 Pm [145] | samarium 62 Sm 150.36 | europium 63 Eu 151.96 | gadolinium 64 Gd 157.25 | terbium 65 Tb 158.93 | dysprosium 66 Dy 162.50 | holmium 67 Ho 164.93 | erbium 68 Er 167.26 | thulium 69 Tm 168.93 | ytterbium 70 Yb 173.04 |
| actinium 89 Ac [227] | thorium 90 Th 232.04 | protactinium 91 Pa 231.04 | uranium 92 U 238.03 | neptunium 93 Np [237] | plutonium 94 Pu [244] | americium 95 Am [243] | curium 96 Cm [247] | berkelium 97 Bk [247] | californium 98 Cf [251] | einsteinium 99 Es [252] | fermium 100 Fm [257] | mendelevium 101 Md [258] | nobelium 102 No [259] |

- **fermionic** alkaline-earths have nuclear spin $I > 0$

Electronic level structure of ^{87}Sr (nuclear spin levels not shown)



Brief Motivation and Overview

Unique properties of **fermionic** alkaline-earths:
- metastable optically excited state 3P_0
- in 1S_0 and in 3P_0 , nuclear spin I decoupled from J

Atomic clock experiments

3P_0 —
 1S_0 —

Quantum information with alkaline-earths

Quantum simulation with alkaline-earths

Quantum computing

Rich physics

Insights into condensed matter problems

Quantum Information Processing with Ultracold Alkaline-Earth Atoms

Quantum information with ultracold fermionic alkaline-earths

Proposals:

- Derevianko, Cannon PRA '04
- Hayes, Julienne, Deutsch PRL '07
- Reichenbach, Deutsch PRL '07
- Daley, Boyd, Ye, Zoller PRL '08
- AG, Rey, Daley, Boyd, Ye, Zoller, Lukin PRL '09
- Reichenbach, Julienne, Deutsch PRA '09
- Shibata, Kato, Yamaguchi, Uetake, Takahashi
App. Phys. B '09
- etc...

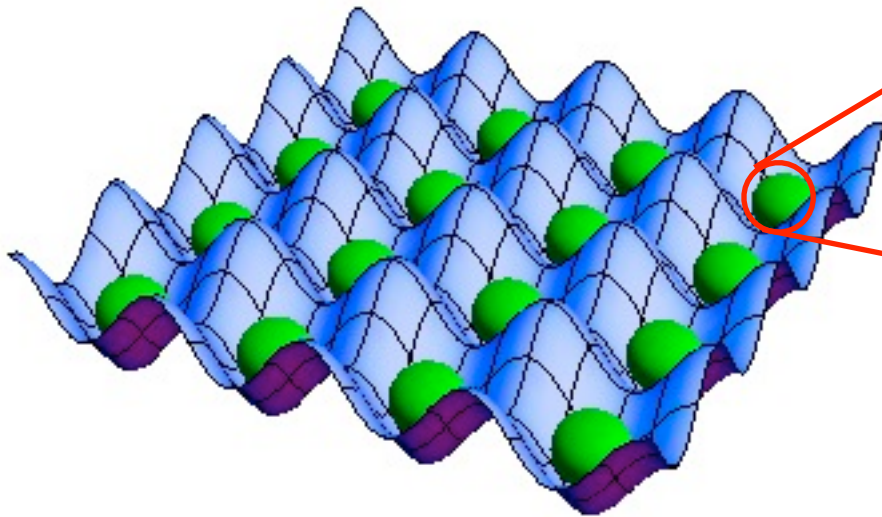
Alkaline-Earth Atoms as Few-Qubit Quantum Registers

The Idea

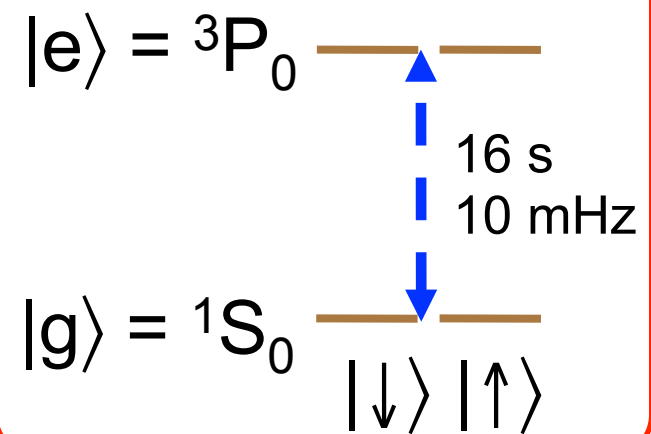
Need for accurate manipulation of large quantum systems:

- quantum computation, precision measurements, etc...

Array of few-qubit quantum registers!



Ex: ^{171}Yb ($I = 1/2$)



Need to:

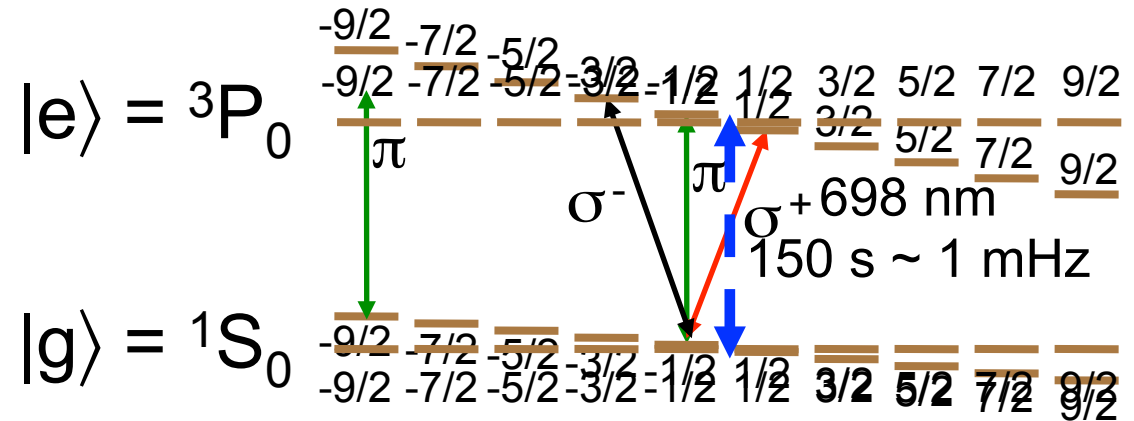
- 1) accurately manipulate single register
- 2) detect $|g\rangle$ - $|e\rangle$ qubit without destroying $|\downarrow\rangle$ - $|\uparrow\rangle$ qubit
- 3) couple registers

Related: Cirac et al.,
Dur & Briegel, Sorensen
& Molmer, Jiang et al.,
Monroe, Saffman,
Hayes et al, Daley et al,
Stock et al, Strauch et
al...

Manipulate an Individual Register: Easy

Ex: ^{87}Sr ($I = 9/2$) [1 electronic + up to 3 nuclear qubits]

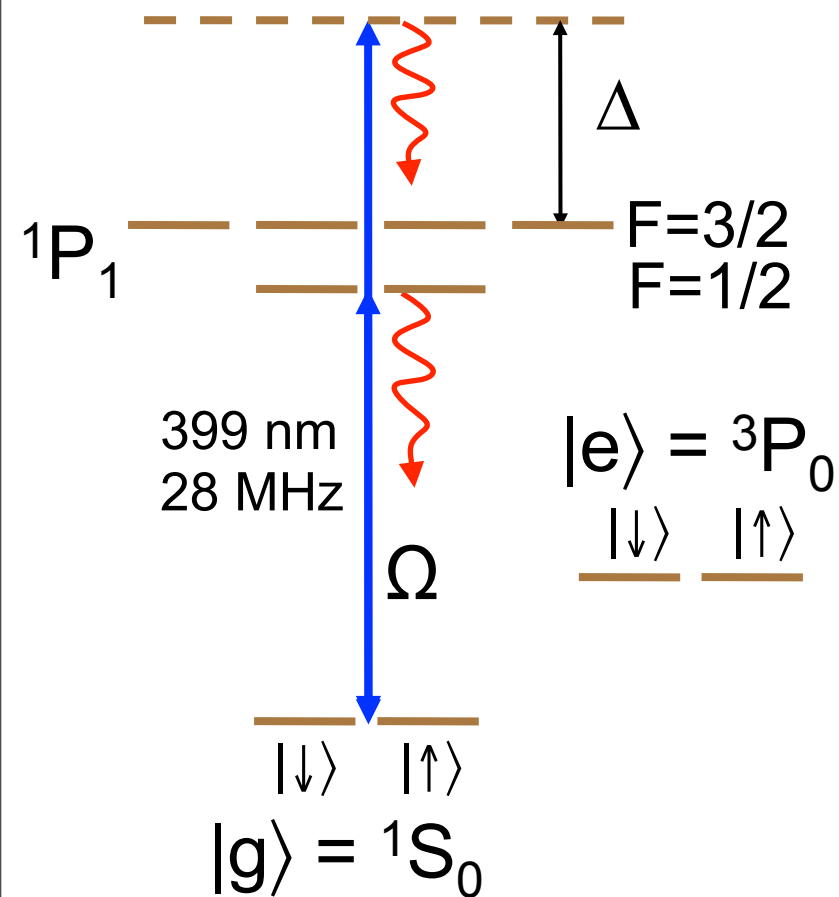
- apply a magnetic field
- different g-factors



- Transitions resolved in experiments [Boyd et al., Science (2006)].

Detect the Electronic Qubit without Erasing the Nuclear Qubits

Ex: ^{171}Yb ($I = 1/2$)



Problem: hyperfine coupling in 1P_1

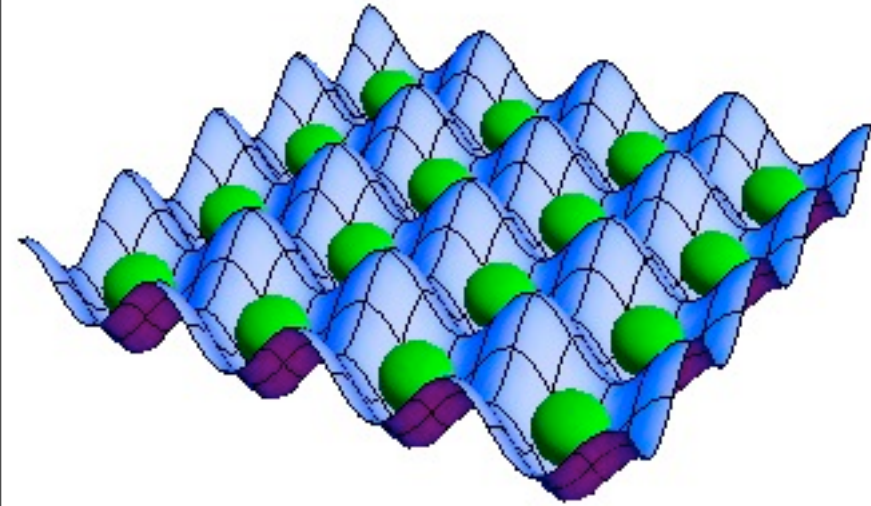
$$\hat{H} = A\hat{\mathbf{I}} \cdot \hat{\mathbf{J}}$$

Solution: Use off-resonant fluorescence
[Childress et al, PRA (2005)]

- destructive interference prevents nuclear spin flips

Related work: Reichenbach & Deutsch, PRL (2007)

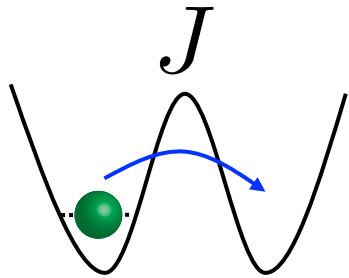
Coupling two registers



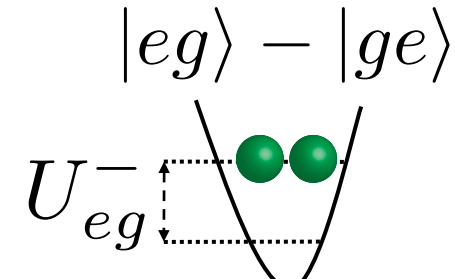
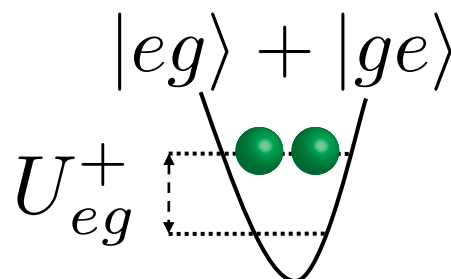
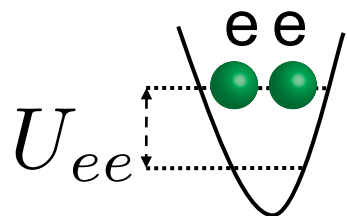
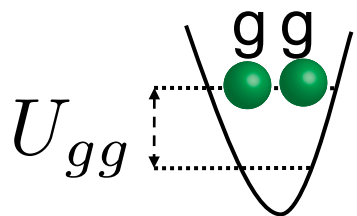
$$|e\rangle = {}^3P_0 \text{ —}$$

$$|g\rangle = {}^1S_0 \text{ —}$$

- Tunneling rate (assume same for g and e):

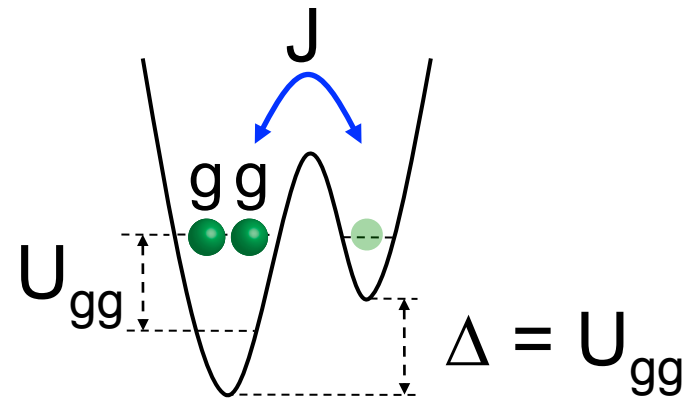
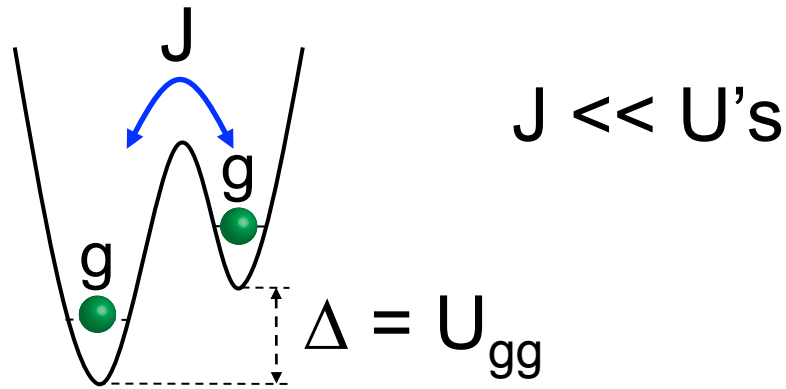


- 4 **different** interaction energies:



Coupling two registers

Ex: ^{171}Yb ($I = 1/2$): spin states \uparrow and \downarrow



- resonantly driven two-level system [Folling et al, Nature (2007), Cheinet et al, PRL (2008)]
- after time $T \propto 1/J$, have a 2π pulse:

$$|g,g\rangle(|\uparrow,\downarrow\rangle - |\downarrow,\uparrow\rangle) \rightarrow -|g,g\rangle(|\uparrow,\downarrow\rangle - |\downarrow,\uparrow\rangle)$$

- repeat for all nuclear Bell states

=> **two-qubit phase gate on the electrons:**

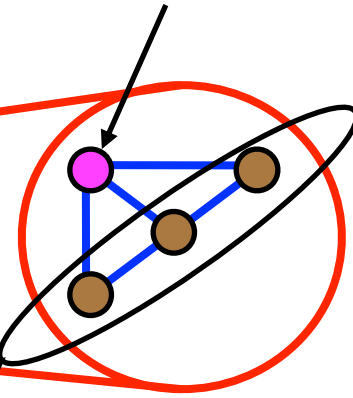
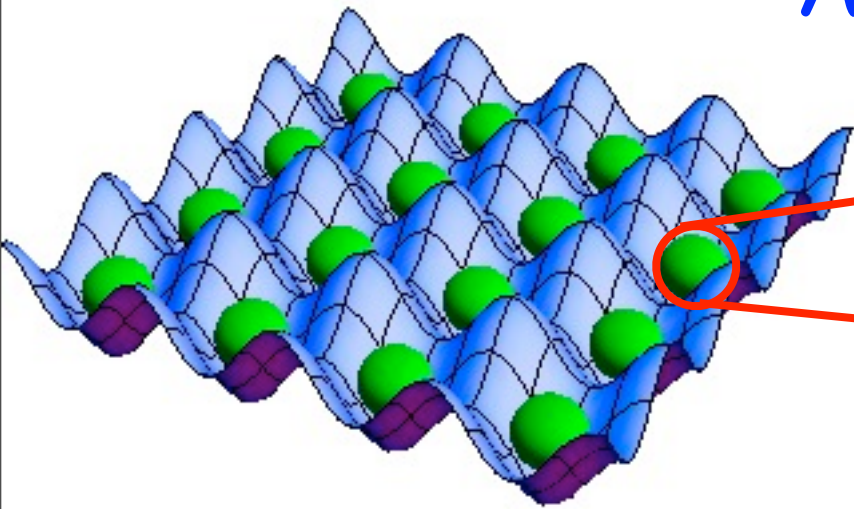
$$|g,g\rangle \rightarrow -|g,g\rangle \quad |g,e\rangle \rightarrow |g,e\rangle \quad |e,g\rangle \rightarrow |e,g\rangle \quad |e,e\rangle \rightarrow |e,e\rangle$$

=> universal manipulation of the full multiregister system

- similarly for $I > 1/2$ Related ideas: Hayes et al, PRL (2007); Daley et al, PRL (2008); Stock et al, PRA (2008); Strauch et al, PRA (2008)

Applications

electronic qubit



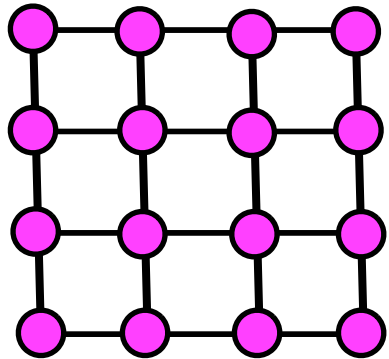
a few (1-3) nuclear qubits

=> circuit-based quantum computation [Jiang et al., PRA (2007)]

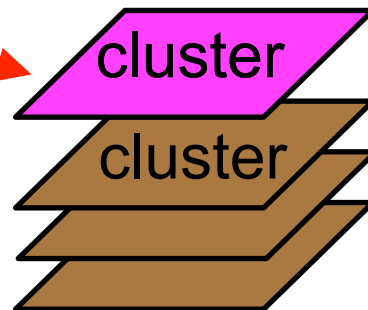
=> generation of high-fidelity entangled many-body states

[Dur & Briegel, PRL (2003)]

- **cluster states** (for measurement-based quantum computation)
- GHZ states (for precision measurements)



$$|+\rangle|+\rangle \cdots |+\rangle$$
$$|+\rangle = |g\rangle + |e\rangle$$



entanglement
pumping/purification

Briegel, Dur, Cirac, Zoller, PRL (1998)

Dur & Briegel, PRL (2003)

Aschauer, Dur, Briegel, PRA (2005)

Details: **PRL 102, 110503 (2009)**

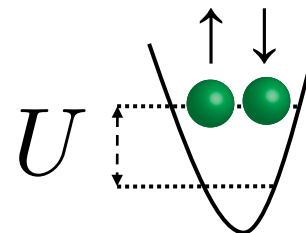
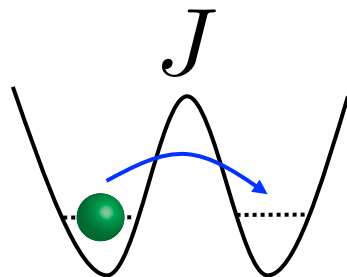
Quantum Simulation with Alkaline-Earth Atoms in Optical Lattices: Two-Orbital $SU(N)$ Magnetism

Outline

- Two orbitals and $SU(N)$ symmetry
- Symmetries of the Hamiltonian in more detail
- Four examples:
 - Examples 1 & 2: $SU(N)$ symmetry
 - Examples 3 & 4: two orbitals

Reminder: the usual Hubbard model

$$H = -J \sum_{\langle j,i \rangle, \sigma} c_{i\sigma}^\dagger c_{j\sigma} + U \sum_j n_{j\uparrow} n_{j\downarrow}$$



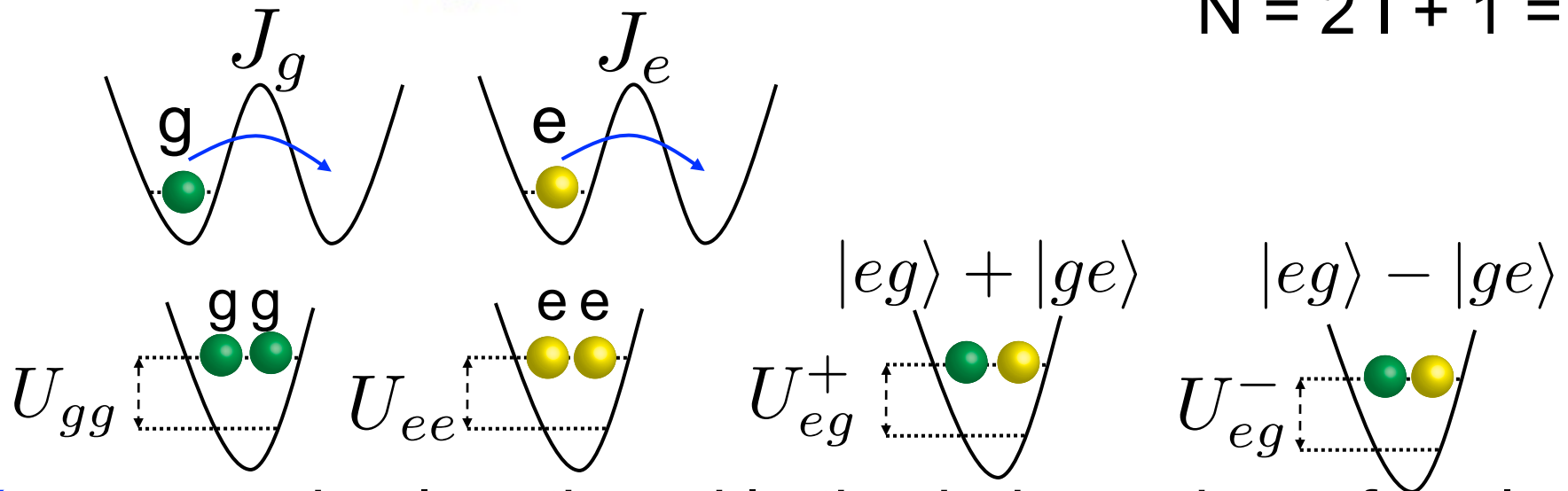
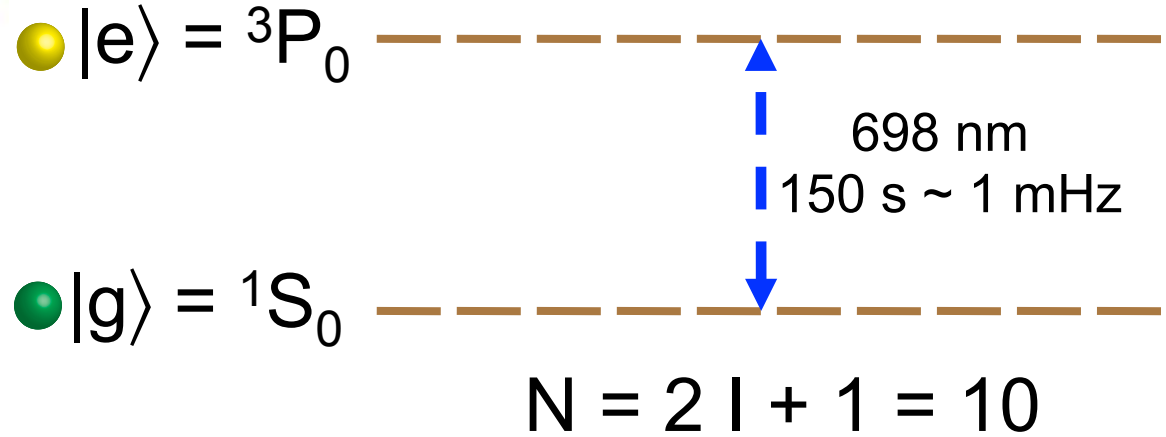
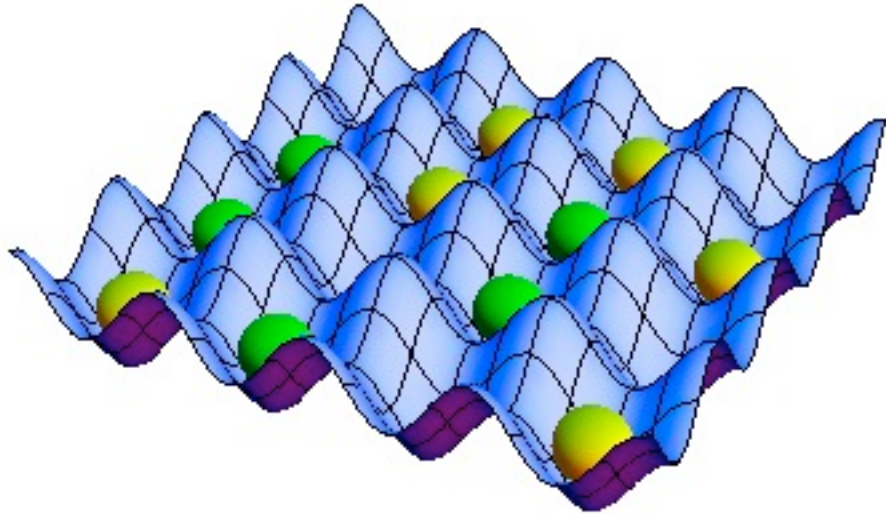
$c_{j\sigma}^\dagger$ - fermion on site j , with spin $\sigma = \uparrow, \downarrow$

$$n_{j\sigma} = c_{j\sigma}^\dagger c_{j\sigma}$$

- high-temperature superconductivity
- cold-atom realizations: Esslinger, Bloch, ...

Two orbitals & $SU(N)$ symmetry

Ex: ^{87}Sr ($I = 9/2$)

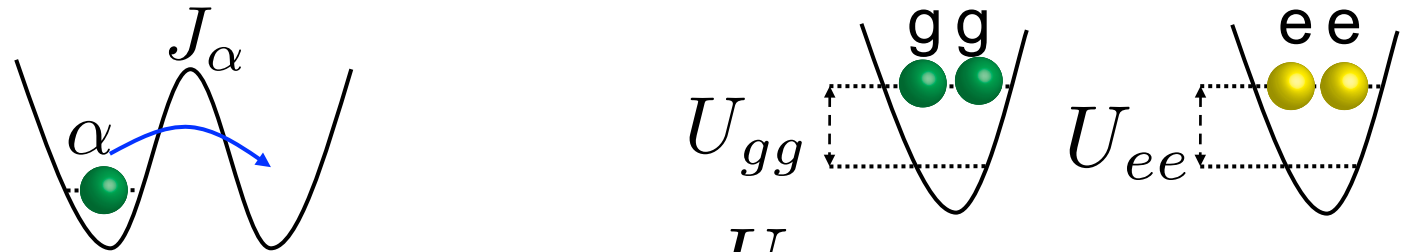


Key: scattering length and lattice independent of nuclear spin (aside from fermionic statistics)

$\Rightarrow SU(N)$ symmetry

See also Cazalilla, Ho, Ueda, New J. Phys. (2009): $SU(6)$ symmetry in ^{173}Yb

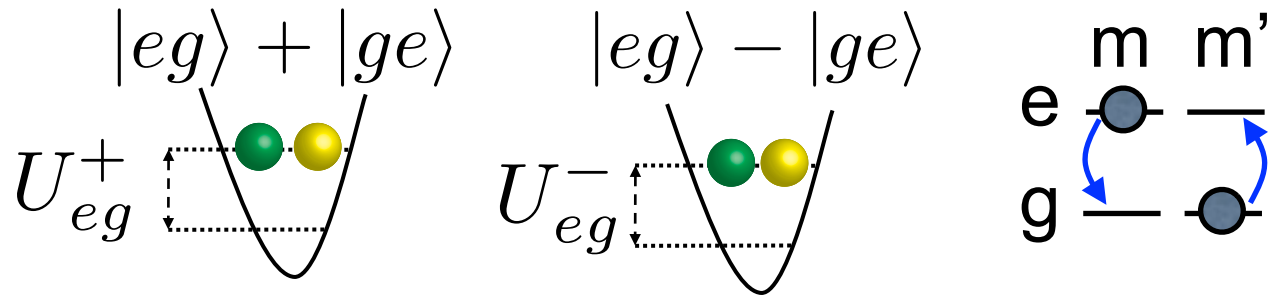
Two-orbital SU(N)-symmetric single-band Hubbard model



$$H = - \sum_{\langle j,i \rangle \alpha, m} J_\alpha c_{i\alpha m}^\dagger c_{j\alpha m} + \sum_{j, \alpha} \frac{U_{\alpha\alpha}}{2} n_{j\alpha} (n_{j\alpha} - 1) + V \sum_j n_{je} n_{jg} + V_{ex} \sum_{j, m, m'} c_{jgm}^\dagger c_{jem'}^\dagger c_{jgm'} c_{jem}$$

$$V = (U_{eg}^+ + U_{eg}^-) / 2$$

$$V_{ex} = (U_{eg}^+ - U_{eg}^-) / 2$$



$c_{j\alpha m}^\dagger$ - site j , electronic state α (g or e), nuclear spin m ($-I \dots I$)

$$n_{j\alpha} = \sum_m n_{j\alpha m} = \sum_m c_{j\alpha m}^\dagger c_{j\alpha m}$$

Symmetries of the Hamiltonian

$$\begin{aligned}
 H = & - \sum_{\langle j,i \rangle \alpha, m} J_\alpha c_{i\alpha m}^\dagger c_{j\alpha m} + \sum_{j, \alpha} \frac{U_{\alpha\alpha}}{2} n_{j\alpha} (n_{j\alpha} - 1) \\
 & + V \sum_j n_{je} n_{jg} + V_{ex} \sum_{j, m, m'} c_{jgm}^\dagger c_{jem'}^\dagger c_{jgm'} c_{jem}
 \end{aligned}$$

SU(2) pseudospin algebra

$$T^x = \frac{1}{2} \sum_{jm} (c_{jem}^\dagger c_{jgm} + c_{jgm}^\dagger c_{jem})$$

$$T^y = \frac{1}{2} \sum_{jm} (-i c_{jem}^\dagger c_{jgm} + i c_{jgm}^\dagger c_{jem}),$$

$$T^z = \frac{1}{2} \sum_j (n_{je} - n_{jg}),$$

$$[T^x, T^y] = iT^z \text{ etc...}$$

$$[T^z, H] = 0$$

U(1) x SU(N)

SU(N=2I+1) algebra

generating nuclear spin rotations

$$S_n^m = \sum_{j, \alpha} c_{j\alpha n}^\dagger c_{j\alpha m}$$

$$[S_n^m, S_q^p] = \delta_{mq} S_n^p - \delta_{pn} S_q^m$$

$$[S_n^m, H] = 0 \quad \forall m, n$$

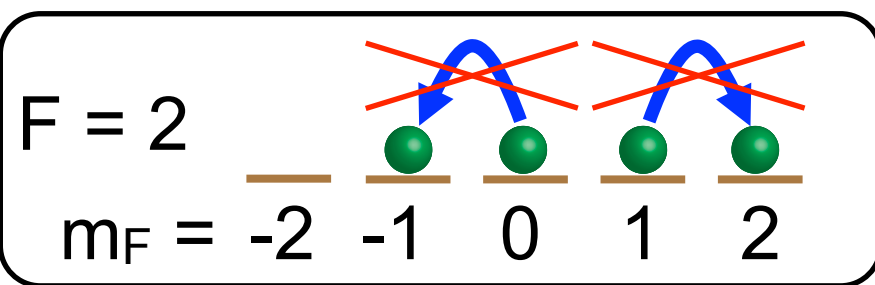
Symmetries of the Hamiltonian

$$H = - \sum_{\langle j,i \rangle \alpha, m} J_\alpha c_{i\alpha m}^\dagger c_{j\alpha m} + \sum_{j, \alpha} \frac{U_{\alpha\alpha}}{2} n_{j\alpha} (n_{j\alpha} - 1) + V \sum_j n_{je} n_{jg} + V_{ex} \sum_{j, m, m'} c_{jgm}^\dagger c_{jem'}^\dagger c_{jgm'} c_{jem}$$

Notice: $S_m^m = \sum_{j\alpha} n_{j\alpha m}$

$$[S_m^m, H] = 0$$

=> if some nuclear Zeeman levels are unoccupied, they will stay unoccupied.

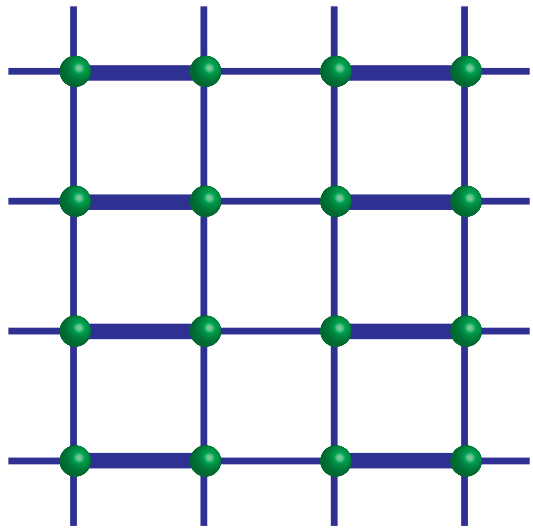


=> I = 9/2 atom (N = 10) can reproduce behavior of all lower I (N < 10) just by choosing initial state

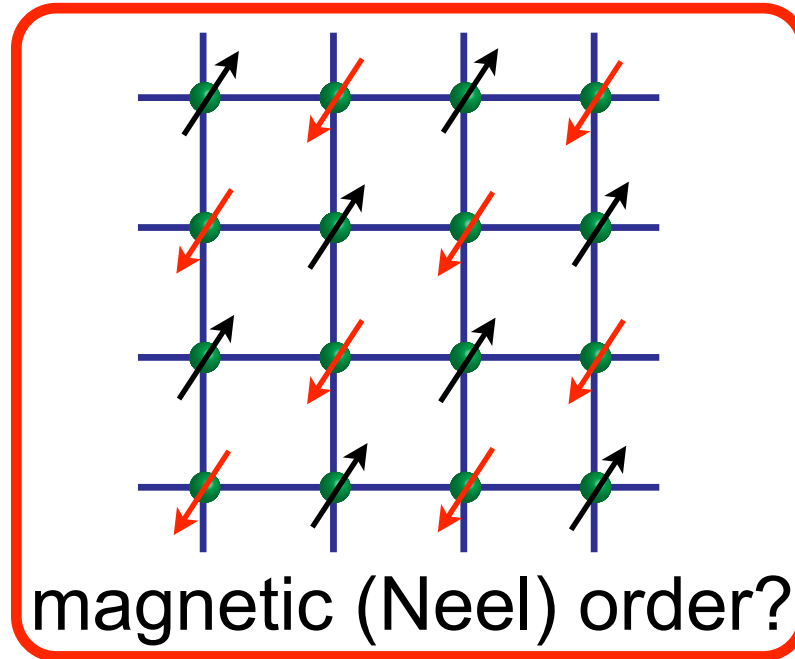
Example 1:

Exotic Valence Bond Solid Phases

$N = 2$ (2 nuclear spin components) one g atom per site



array of singlets?

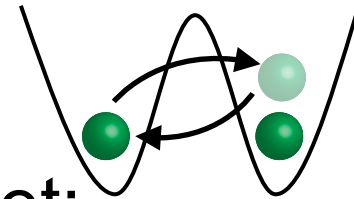


magnetic (Neel) order?

$J_g \ll U_{gg} \Rightarrow$ spin Hamiltonian

SU(2)-symmetric Heisenberg antiferromagnet:

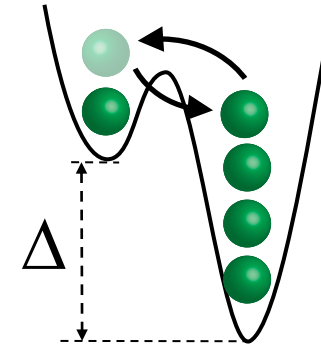
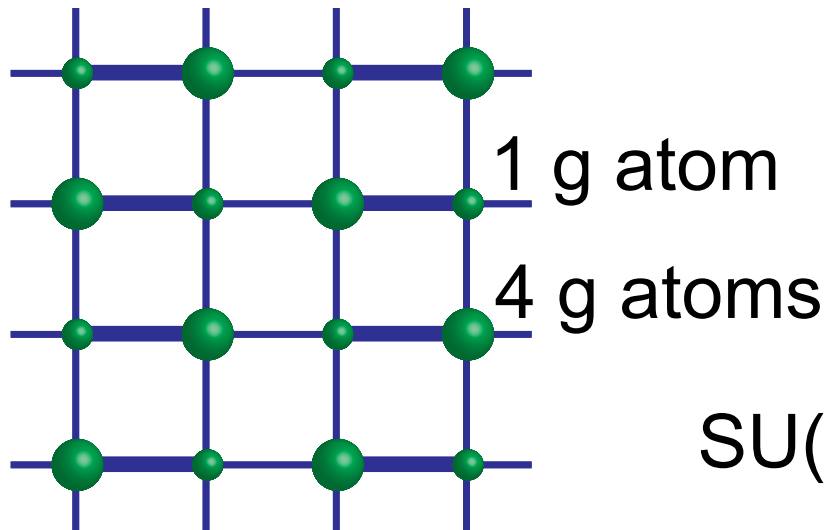
$$H = \frac{4J_g^2}{U_{gg}} \sum_{\langle i,j \rangle} \vec{S}_i \cdot \vec{S}_j$$



Example 1:

Exotic Valence Bond Solid Phases

$N = 5$ (5 nuclear spin components)



$$J_g \ll U_{gg}$$

SU(5)-symmetric spin Hamiltonian

Tune Δ to get anti-ferromagnetic interaction

SU(N) singlets: N atoms, energy $\sim -NJ_g^2/U_{gg}$

=> exotic valence bond solid (VBS) order is expected

=> gapped & broken rotational symmetry

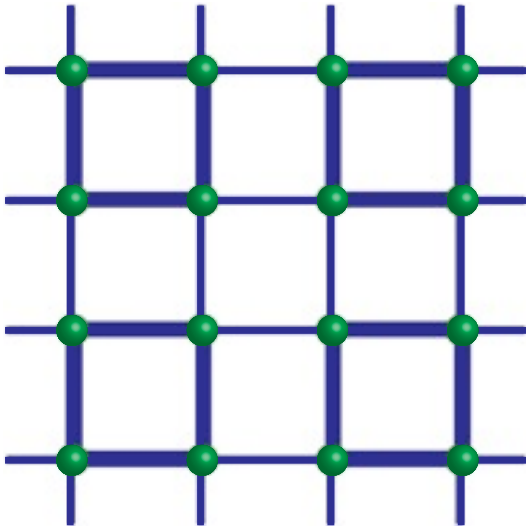
Melting **experiment.**

[Read & Sachdev (1990), Harada et al (2003), Assaad (2005),...]

Example 2:

Exotic Valence Plaquette Solid Phases

$N = 4$ (4 nuclear spin components)



1 g atom per site

[Li et al, PRL (1998); Bossche et al, EPJ B, (2000); Xu & Wu, PRB (2008)]

Similar arrangements => exotic spin liquid phases.

[Hermele, Rey, Gurarie, PRL 103, 135301 (2009)]

Example 3: Kugel-Khomskii model

One atom per site (both g and e allowed)

$$H = \sum_{\langle i,j \rangle} \left[(A + BS_{ij}^2)(T_i^x T_j^x + T_i^y T_j^y) + [C + DS_{ij}^2](T_i^z T_j^z + \frac{1}{4}) + E(1 - S_{ij}^2)(T_i^z + T_j^z) + FS_{ij}^2 \right]$$

~ spin-dependent
XXZ in pseudospin

- competing
orders

$$|mm\rangle \ \& \ |mn\rangle + |nm\rangle \quad N = 2 \Rightarrow \text{triplet}$$

$$S_{ij}^2 = \sum_{mn} S_m^n(i) S_n^m(j) = 1 \text{ if symmetric under } i\text{-}j \text{ exchange}$$

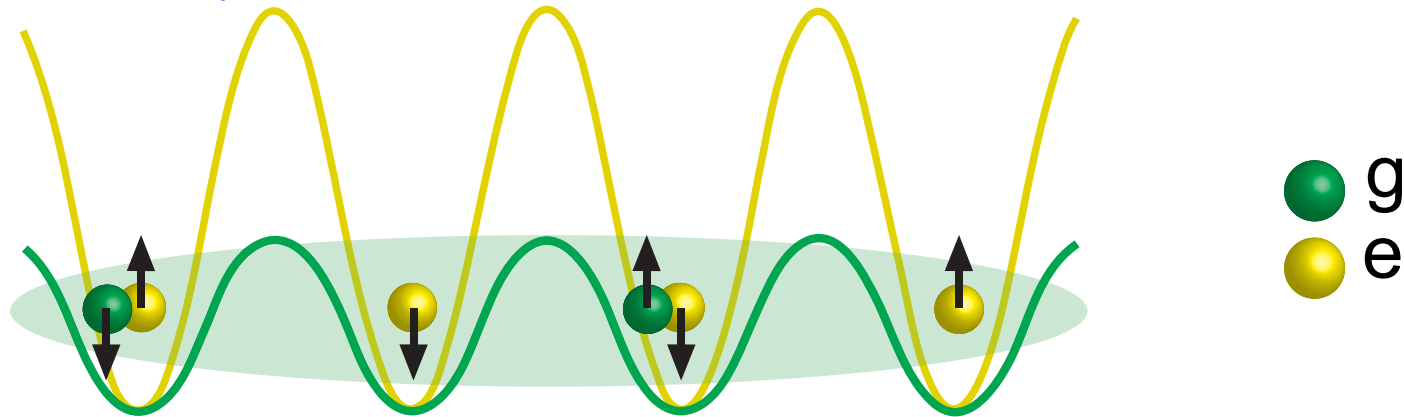
$$= -1 \text{ if antisymmetric under } i\text{-}j \text{ exchange}$$

$$|mn\rangle - |nm\rangle \quad N = 2 \Rightarrow \text{singlet}$$

Generalization to any N of SU(N=2)-symmetric **Kugel-Khomskii model**:
 - used to model spin-orbital interactions in transition metal oxides with perovskite structure [Kugel,Khomskii, JETP (1973); Tokura,Nagaosa, Science (2000)]

Example 4:

SU(N)-symmetric Kondo Lattice Model



$$H_{KL} = - \sum_{\langle j,i \rangle \alpha, m} J_g c_{igm}^\dagger c_{jgm} + V_{ex} \sum_{j, m, m'} c_{jgm}^\dagger c_{jem'}^\dagger c_{jgm'} c_{jem}$$

- N = 2 case is important in the study of
 - colossal magnetoresistance
 - heavy fermion materials
- Temperature requirement is favorable:
e.g. heavy Fermi liquid may be accessed with $k_B T \sim U$
(not J^2/U)

[Proposals to study N = 2 Kondo physics with cold atoms:
Detailed study: Foss-Feig, Hermele, Gurarie, Rev PRA '10; arxiv '10
Duan, Europhys. Lett. '04; Paredes, Tejedor, Cirac, PRA '05]

Conclusions

Alkaline-earth atoms for [quantum information processing](#)

[[PRL 102, 110503 \(2009\)](#)]:

- few-qubit quantum registers
- applications to quantum computing and precision measurements

Alkaline-earth atoms for [quantum simulation](#) [[Nature Phys. 6, 289 \(2010\)](#)]:

- two-orbital $SU(N)$ Hubbard model
 - the Kugel-Khomskii model
 - $SU(N)$ antiferromagnets
 - $SU(N)$ Kondo lattice model
- few-body (skipped) & many-body experiments
- possible insights into strongly-correlated systems in condensed matter

Outlook

- Analyze new models
- Alkali dimers, ions

- Clock shifts:
[PRL 103, 260402 \(2009\)](#)

^{87}Sr ($I = 9/2$)**Degenerate Fermi Gas of ^{87}Sr**

B. J. DeSalvo, M. Yan, P. G. Mickelson, Y. N. Martinez de Escobar, and T. C. Killian

Department of Physics and Astronomy, Rice University, Houston, Texas, 77251, USA

(Received 5 May 2010; published 13 July 2010)

PHYSICAL REVIEW A **82**, 011608(R) (2010)**Double-degenerate Bose-Fermi mixture of strontium**Meng Khoon Tey,¹ Simon Stellmer,^{1,2} Rudolf Grimm,^{1,2} and Florian Schreck¹¹*Institut für Quantenoptik und Quanteninformation, Österreichische Akademie der Wissenschaften, 6020 Innsbruck, Austria*²*Institut für Experimentalphysik und Zentrum für Quantenphysik, Universität Innsbruck, 6020 Innsbruck, Austria*

(Received 6 June 2010; published 20 July 2010)

 ^{171}Yb ($I = 1/2$) and ^{173}Yb ($I = 5/2$) Selected for a Viewpoint in *Physics*

PRL 105, 190401 (2010)

PHYSICAL REVIEW LETTERS

week ending
5 NOVEMBER 2010**Realization of a $\text{SU}(2) \times \text{SU}(6)$ System of Fermions in a Cold Atomic Gas**Shintaro Taie,^{1,*} Yosuke Takasu,¹ Seiji Sugawa,¹ Rekishu Yamazaki,^{1,2} Takuya Tsujimoto,¹
Ryo Murakami,¹ and Yoshiro Takahashi^{1,2}¹*Department of Physics, Graduate School of Science, Kyoto University, Japan 606-8502*²*CREST, JST, 4-1-8 Honcho Kawaguchi, Saitama 332-0012, Japan*

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Conclusions

Alkaline-earth atoms for quantum information processing

[PRL 102, 110503 (2009)]:

- few-qubit quantum registers
- applications to quantum computing and precision measurements

Alkaline-earth atoms for quantum simulation [Nature Phys. 6, 289 (2010)]:

- two-orbital
 - the Kugelblitz model
 - $SU(N)$ Kondo lattice model
 - $SU(N)$ Kondo lattice model
- few-body (skipped) & many-body experiments
- possible insights into strongly-correlated systems in condensed matter

Thank you

Outlook

- Analyze new models
- Alkali dimers, ions

- Clock shifts:
PRL 103, 260402 (2009)