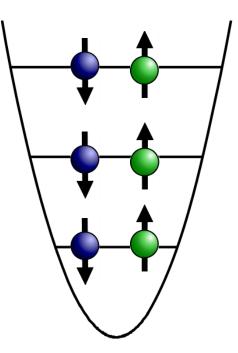


## One, two, three, many Creating quantum systems one atom at a time

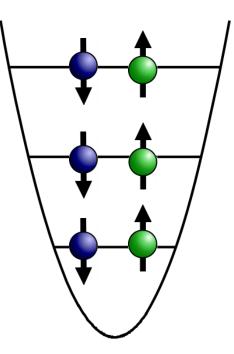


#### Selim Jochim, Universität Heidelberg





## One, two, three, many Creating quantum systems one atom at a time

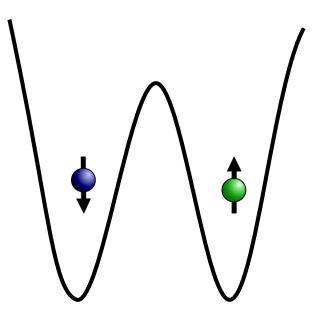


#### Selim Jochim, Universität Heidelberg





## One, two, three, many Creating quantum systems one atom at a time



Selim Jochim, Universität Heidelberg



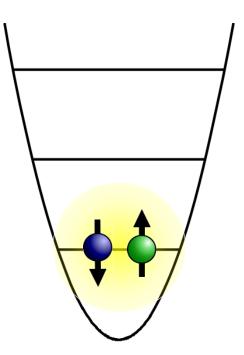


PHYSIKALISCHE

ΙΝSΤΙ



interacting singlet

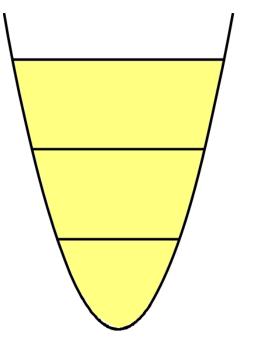


Ground state of the Helium atom:

No analytic solution available, we learn how to apply powerful numerical techniques: Hartree Fock method.







Define quantities like the Fermi energy, density, pressure ....

... apply local density approximation ...

But when are such approximations justified?

This is an ancient problem!



#### **Sorites Paradox**

#### How many grains make a heap?

- 1 grain of sand does not make a heap.
- If 1 grain does not make a heap then 2 grains of sand do not.
- If 2 grains do not make a heap then 3 grains do not.

 If 9,999 grains do not make a heap then 10,000 do not.
 From Stanford Encyclopedia of Philosophy: <u>http://plato.stanford.edu/entries/sorites-paradox/</u>

http://commons.wikimedia.org/wiki/File%3ASossusvlei\_Dune\_Namib\_Desert\_Namibia\_Luca\_Galuzzi\_2004.JPG





Bose Einstein condensates of large samples of atoms: Macroscopic wave function: Number of particles is so large that a constant density of atoms is observed in experiments:

Measure:  $n(\boldsymbol{r}) = \langle \widehat{\Psi}^{\dagger}(\boldsymbol{r}) \widehat{\Psi}(\boldsymbol{r}) \rangle$ 

PHYSIKALISCHE

NS

http://jila.colorado.edu/bec/images/bec.png

Removing one single atom does not make a difference!



### **Reduce the complexity** of a system as much as possible

### until only the essential parts remain!

In most physical systems:

## **Range of interaction**

significantly complicates the description





## The interactions between ultracold atoms can be effectively pointlike (contact interaction)

van der Waals interaction: range of  $r_{vdW} \sim 1$ nm

In the experiments we have:

- extremely low density (interparticle spacing ~ 1µm)
- extremely low momentum, such that  $\lambda_{dB} = \frac{h}{\sqrt{2\pi m kT}} \gg r_{vdW}$





• extremely low momentum, such that  $\lambda_{dB} = \frac{h}{\sqrt{2\pi m kT}} \gg r_{vdW}$ 

(This is the opposite limit desired in collision experiments: shorter wavelength enhances resolution)

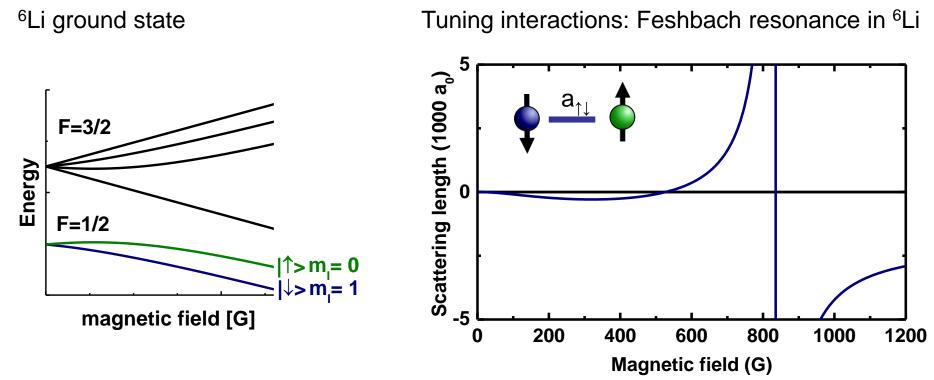
Here:

- If  $\lambda_{dB}$  is sufficiently large, all the information about internal structure of the atom is hidden in a single quantity, **the scattering length** *a*
- We can even tune the scattering length to any desired value by simply applying a magnetic field (**Feshbach resonances**).



#### The <sup>6</sup>Li atom





S=1/2, I=1

 $\rightarrow$  half-integer total angular momentum  $\rightarrow$  <sup>6</sup>Li is a fermion

**NO** interaction between identical particles

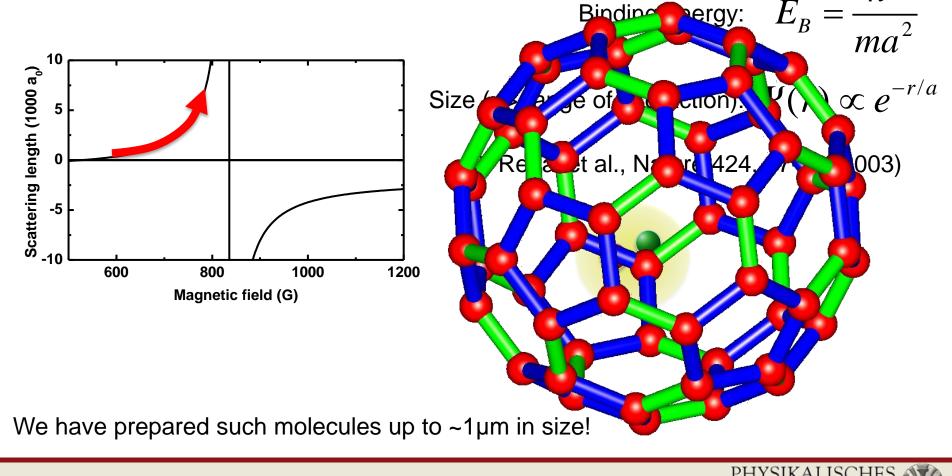
PHYSIKALISCHES INSTITUT

## **Tunability of ultracold systems**



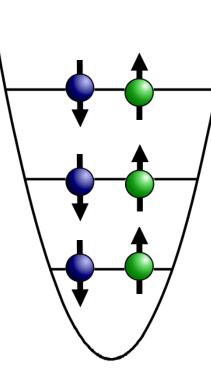
INS

Feshbach resonance: Magneticfield dependence of s-wave scattering length Two-body system: Tune the binding energy of a weakly bound molecule:



G. Zürn et al., PRL 110, 135301 (2013)















How do we prepare our samples?

• How many particles do we need to form a heap?

• Controlling the motion of two particles in a double well

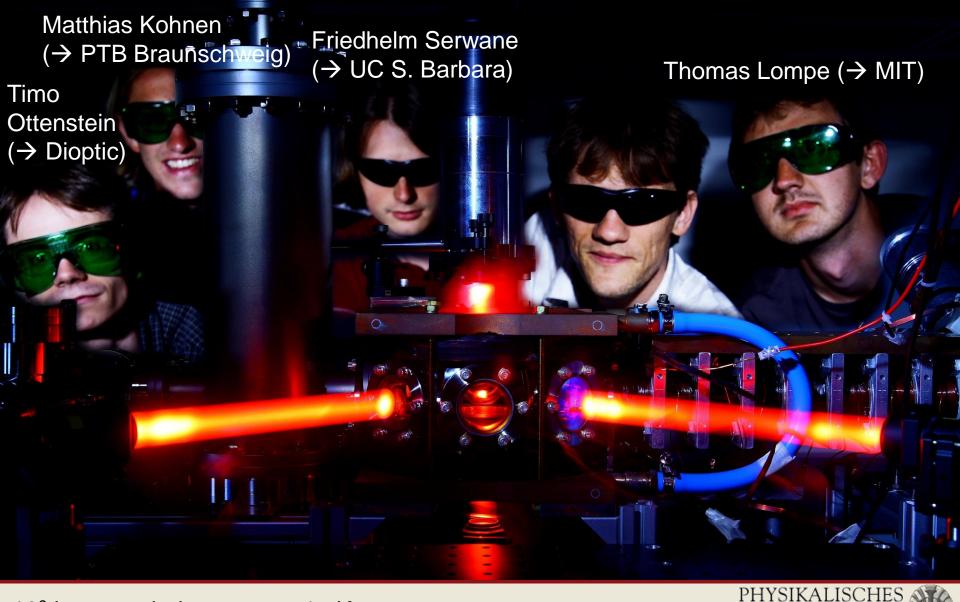


#### A picture from the lab ...



ΙΝSΤ





10<sup>9</sup> laser cooled atoms at ~1mK





We need to isolate the atoms from the environment:



... here we use the focus of a laser beam:



**Optical dipole trap** depth:  $U \propto I(r)$ 

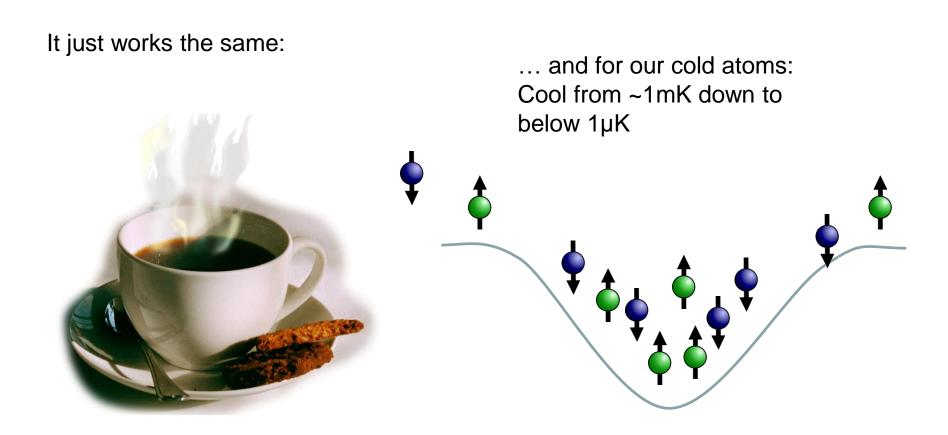
This might still work for liquid nitrogen ....



#### **Evaporative cooling**







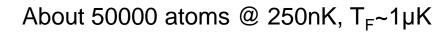
For our cup of coffee ...

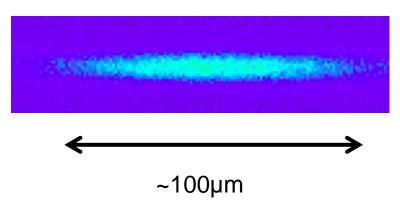
Just reduce the trap depth, i.e. laser power

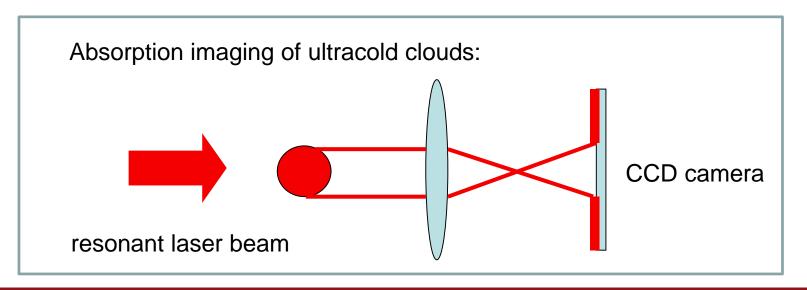








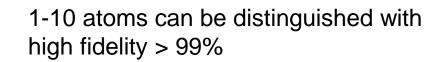


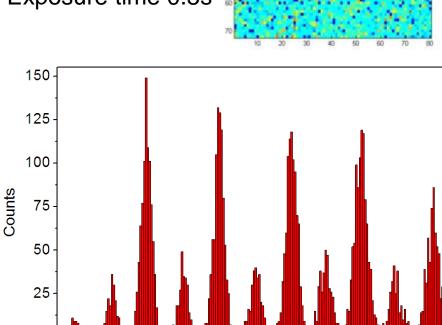


PHYSIKALISCHES INSTITUT

#### Single atom detection

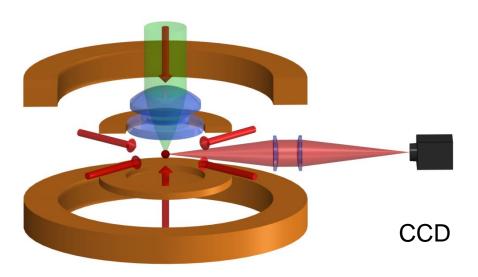
one atom in a MOT 1/e-lifetime: 250s Exposure time 0.5s

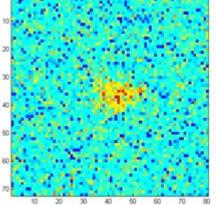




Fluorescence normalized to atom number







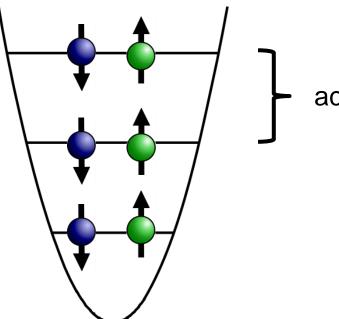


#### Towards a finite gas ...

Universität Heidelberg



The challenge:



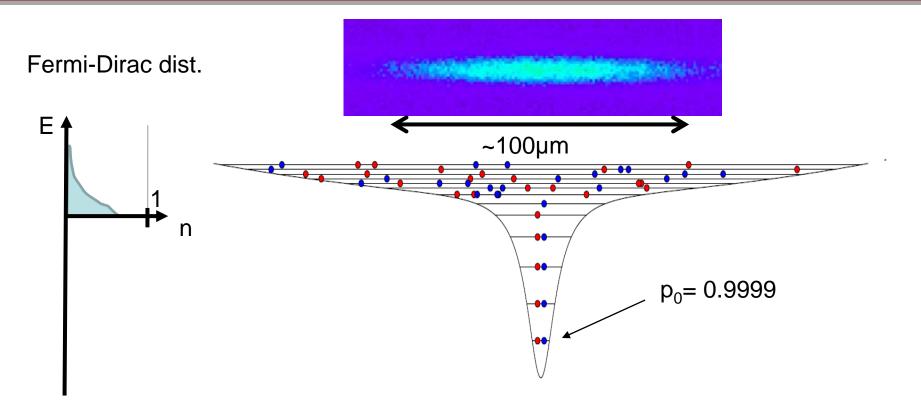
- achieve  $\hbar\omega\gg kT$ 



### Creating a finite gas of fermions







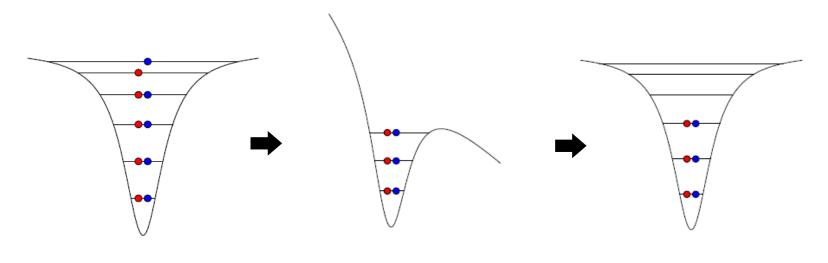
- 2-component mixture in reservoir
- superimpose microtrap (~1.8 µm waist)







switch off reservoir



+ magnetic field gradient in axial direction



#### Spilling the atoms ....

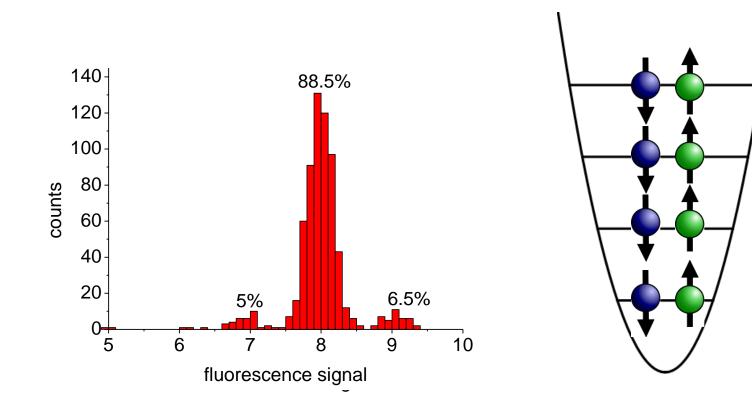
Universität Heidelberg

PHYSIKALISCHES

Т

INS





- We can control the atom number with exceptional precision!
- Note aspect ratio 1:10: 1-D situation
- So far: Interactions tuned to zero ...

F. Serwane et al., Science **332**, 336 (2011)



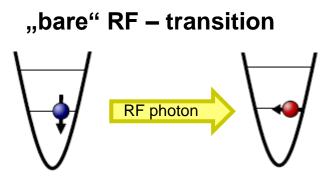
# Let's study the interacting system!



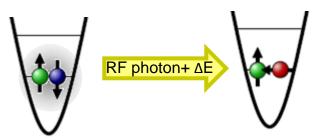




#### Radio Frequency spectroscopy



**RF** – transition with interaction

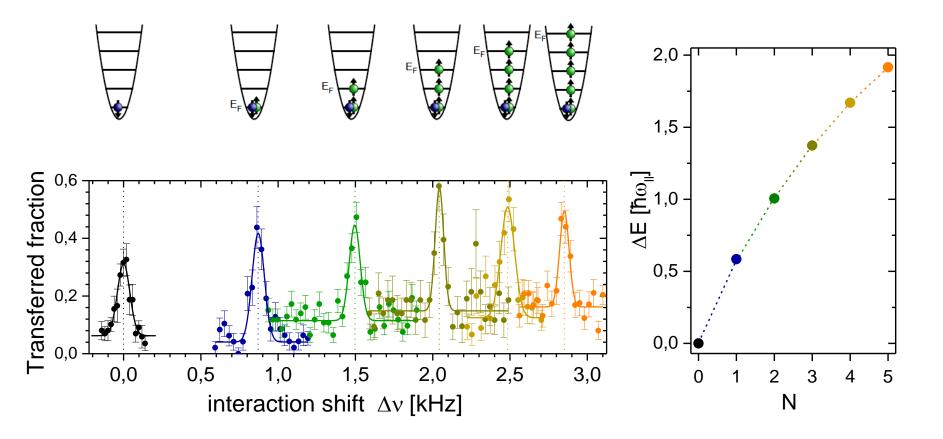


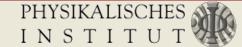






vary the number of majority particles:

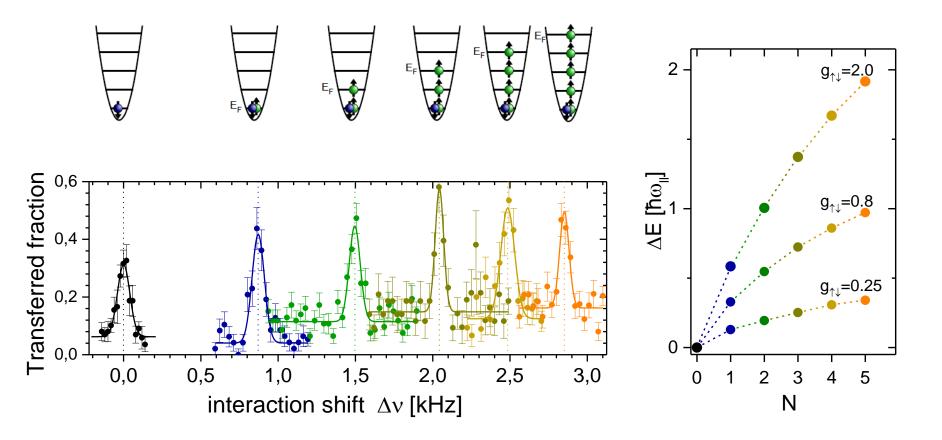




Universität Heidelberg



vary the number of majority particles:





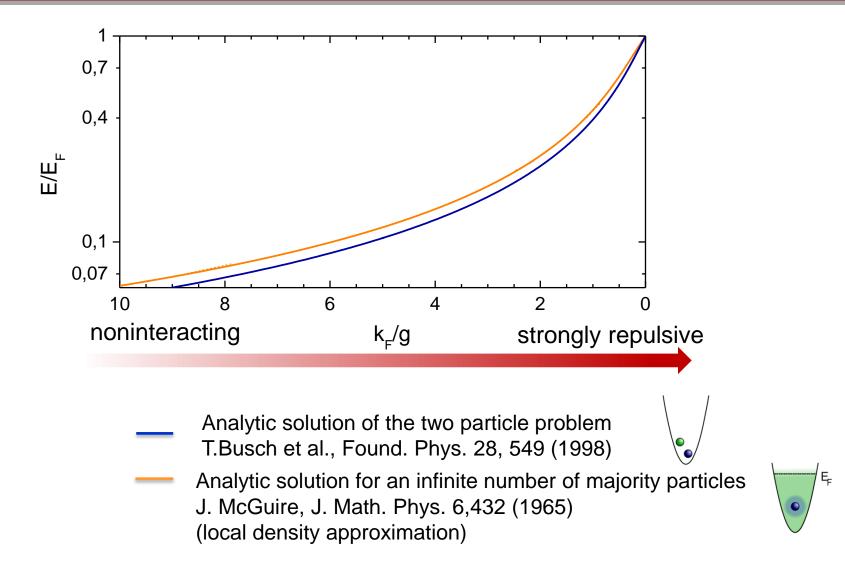
#### Interaction energy in dimensionless units



PHYSIKALISCHES

ΙΝSΤ



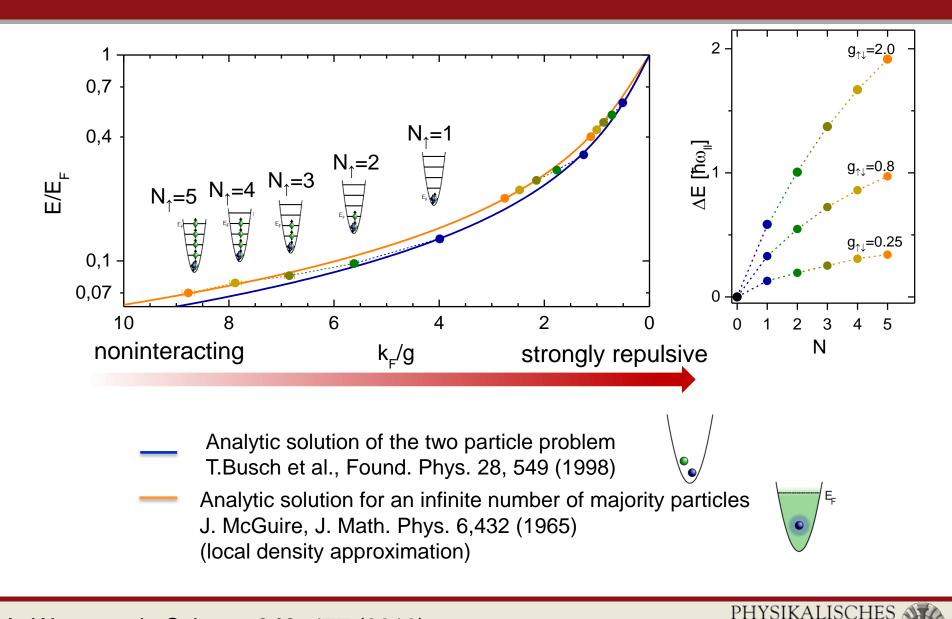


#### Interaction energy in dimensionless units



INS



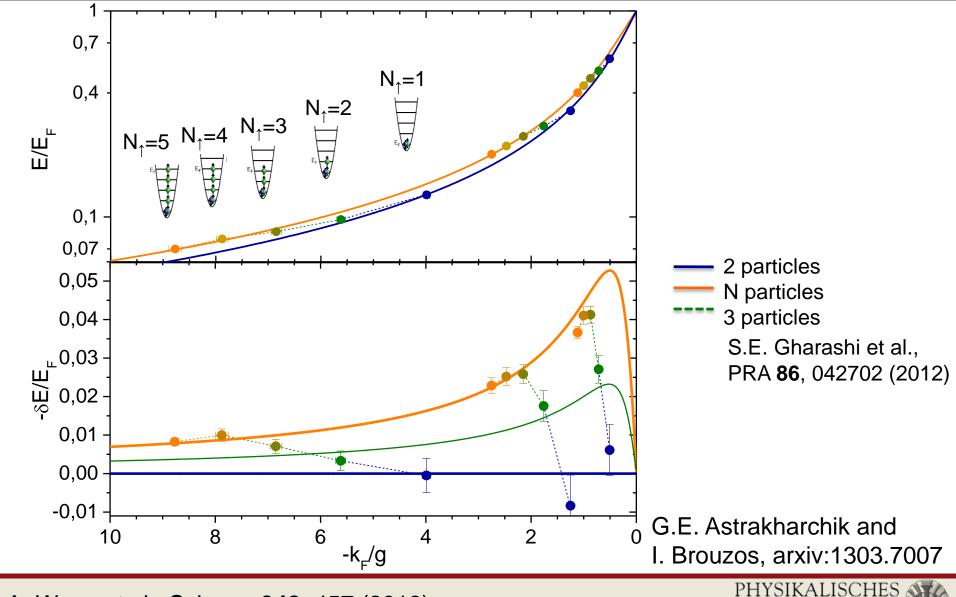


#### Interaction energy in dimensionless units

Universität Heidelberg

ΙΝSΤΙ



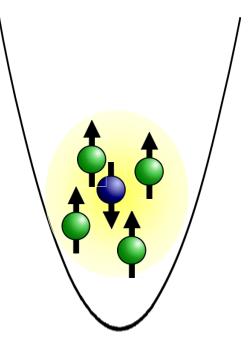


A. Wenz et al., Science 342, 457 (2013)





... with very few particles (in a one-dimensional system)



Interesting things to look at:

- Polaron physics in various dimensions
- The Kondo problem
- Anderson's orthogonality catastrophe

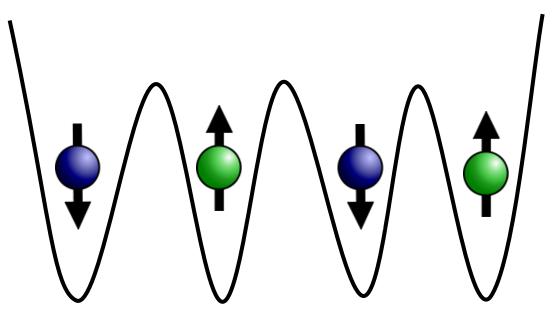


#### A. Wenz et al., Science 342, 457 (2013)





..... with similar fidelity and control?



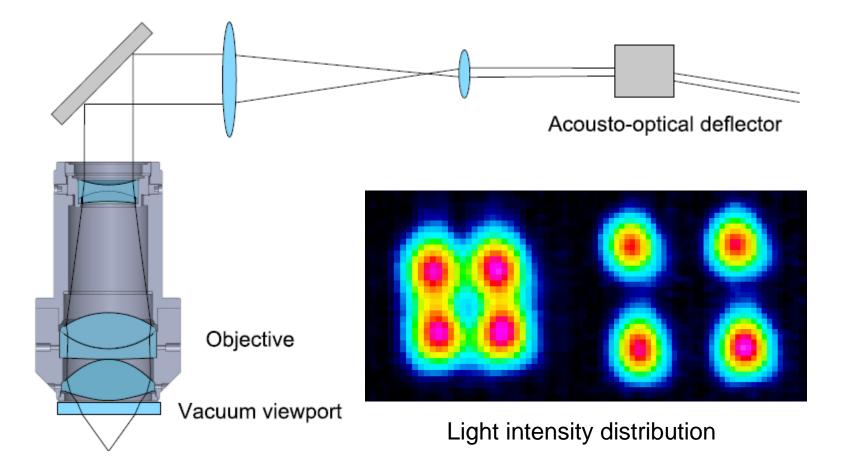
Basic building blocks of matter!



#### The setup









#### A tunable double well

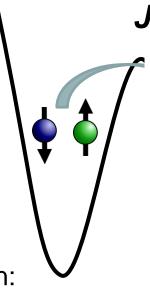


PHYSIKALISCHES

Τ

INSTI





initial spatial wave function:

$$|\Psi(t=0)
angle = |L
angle_1 |L
angle_2 = |LL
angle$$

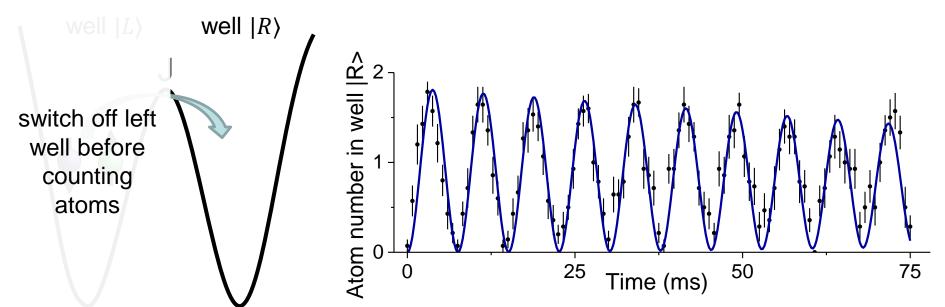
spin wave function (stationary)

$$|\chi\rangle = \frac{1}{\sqrt{2}} (|\uparrow\rangle_1 |\downarrow\rangle_2 - |\downarrow\rangle_1 |\uparrow\rangle_2) = \frac{1}{\sqrt{2}} (|\uparrow\downarrow\rangle - |\downarrow\uparrow\rangle)$$

Universität Heidelberg







$$\begin{split} |\Psi(t)\rangle &= |\psi(t)\rangle_1 \, |\psi(t)\rangle_2 \\ |\psi(t)\rangle_1 &= \frac{1}{2}((|L\rangle_1 + |R\rangle_1) + (|L\rangle_1 - |R\rangle_1)e^{-i\Delta Et/\hbar}) \end{split}$$

PHYSIKALISCHES INSTITUT

#### Two interacting atoms

Universität Heidelberg

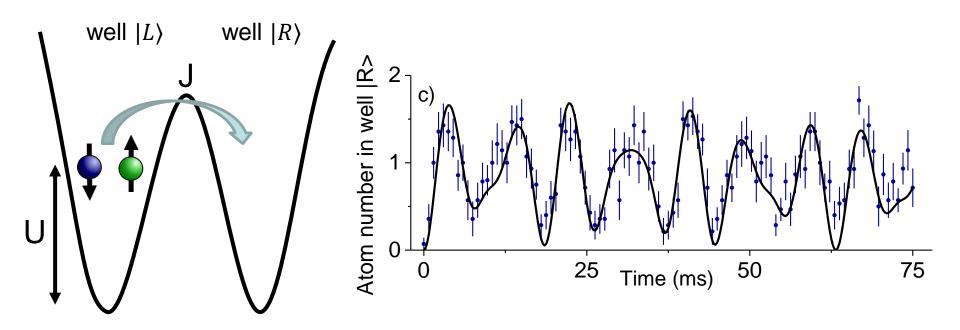
PHYSIKALISCHES

Т

INS



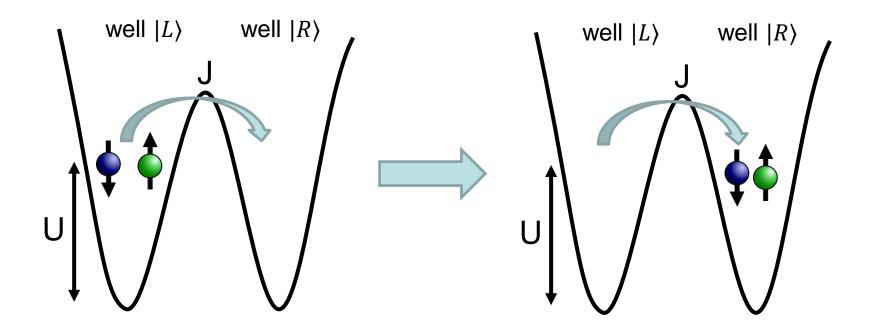
Interaction leads to entanglement:



 $|\Psi(t)
angle 
eq |\psi(t)
angle_1 |\psi(t)
angle_2$ 





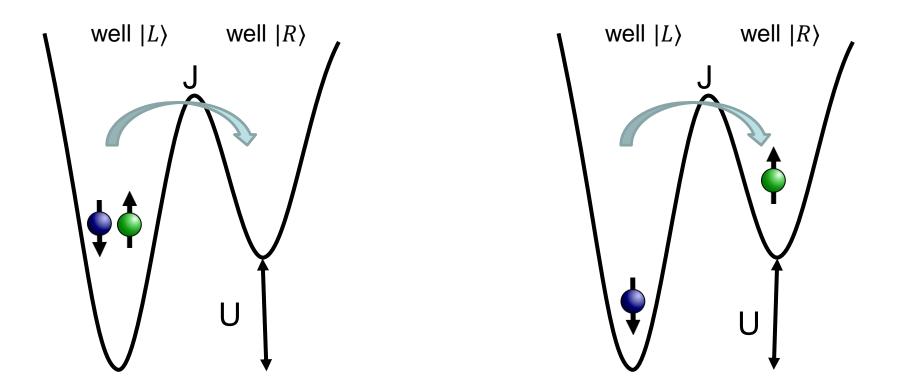


In a balanced double well, they can only tunnel together!





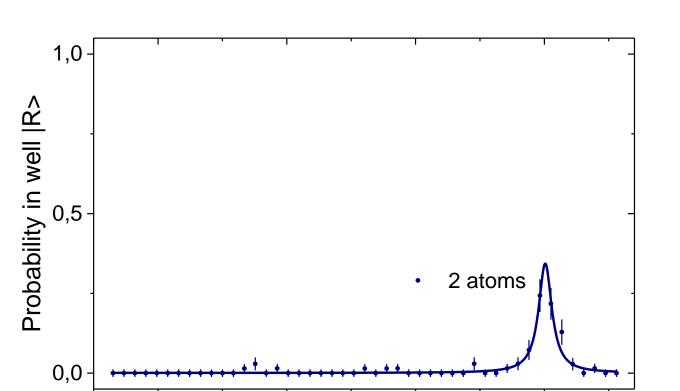




We can compensate for the interaction energy by applying a tilt!







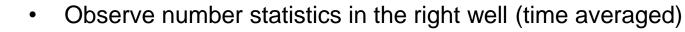
tilt [J]

-5

0

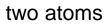
B=740G

-10



Two strongly interacting atoms

-15

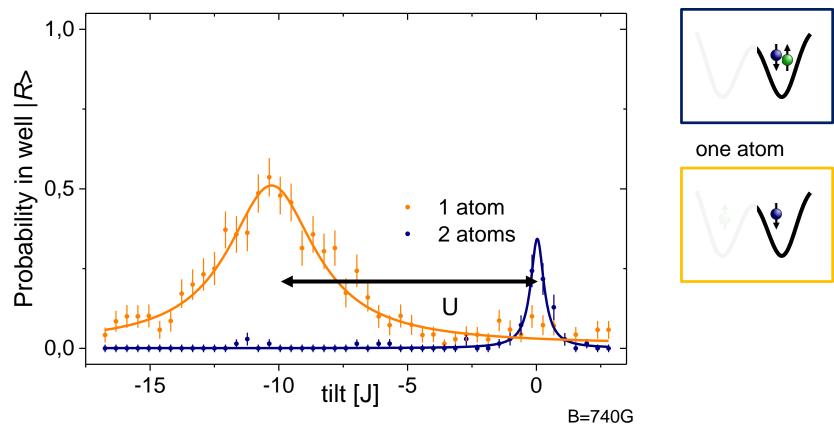


Universität

Heidelberg



• Observe number statistics in the right well (time averaged)



two atoms

PHYSIKALISCHES

ΙΝ S Τ

**U**NIVERSITÄT

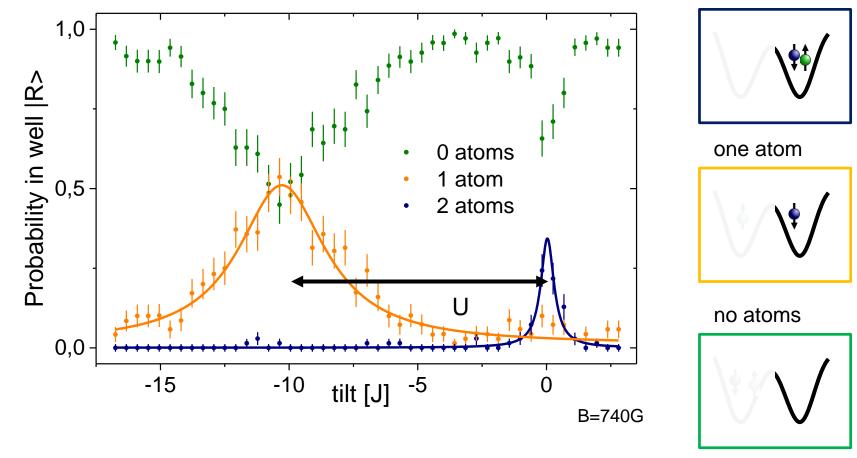
Heidelberg



#### PHYSIKALISCHES INSTITUT



• Observe number statistics in the right well (time averaged)



two atoms

Universität

Heidelberg

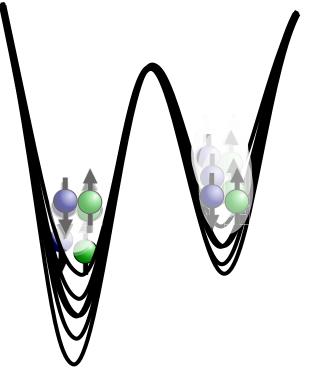


### **Preparing stationary states**





• If we ramp on the second well slowly enough, the system will remain in its ground state:

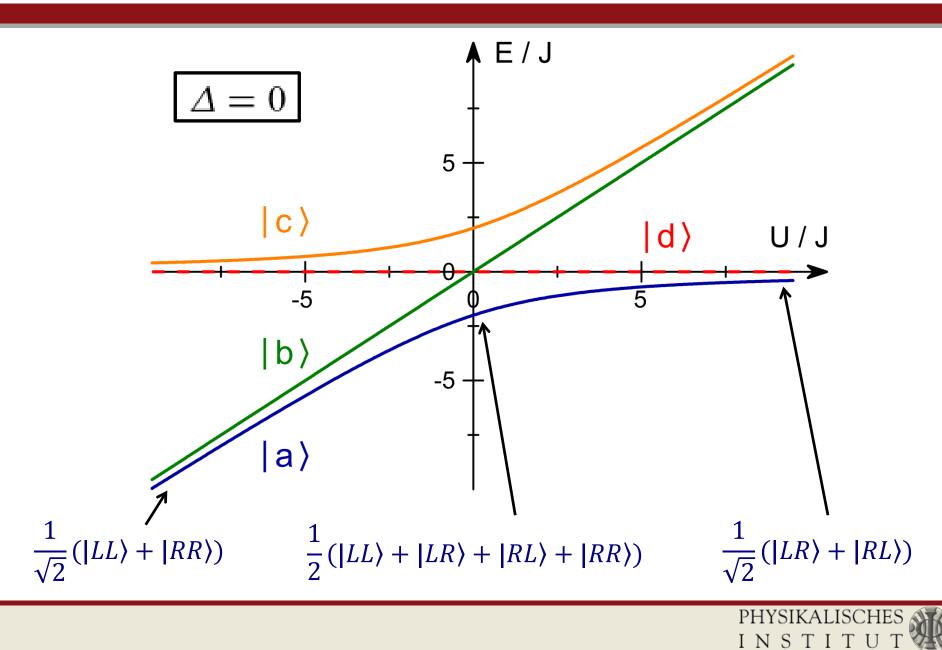




#### **Eigenstates of a symmetric DW**

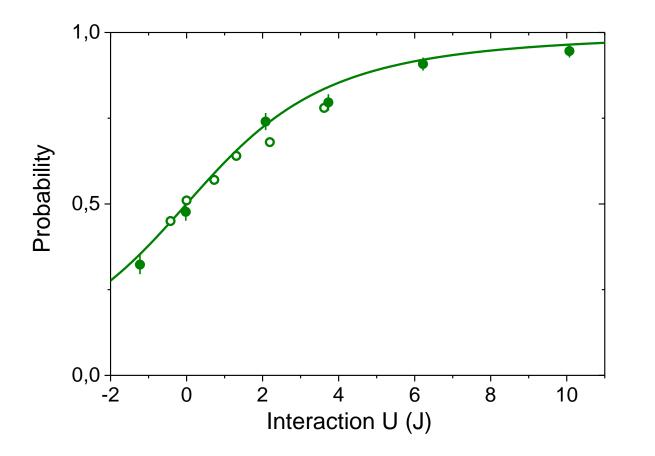




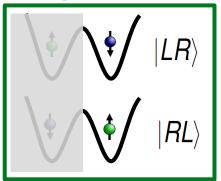


## **Preparing stationary states**

• Number statistics for the balanced case depending on the interaction strength:



#### Single occupancy

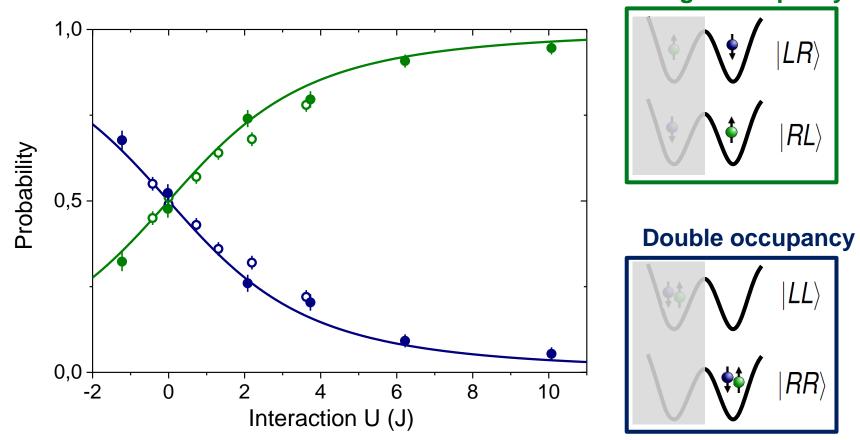


PHYSIKALISCHES



## **Preparing stationary states**

Number statistics for the balanced case depending on the interaction strength:



#### Single occupancy

Universität

Heidelberg



PHYSIKALISCHES INSTITUT

#### **Measuring energies**

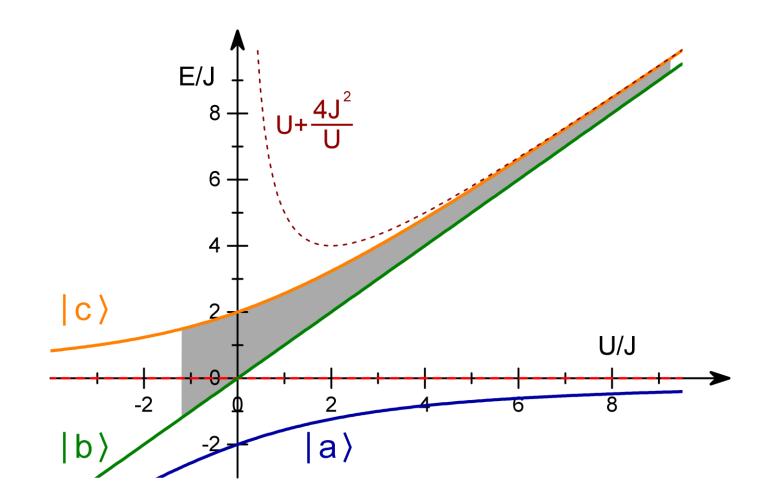
Universität Heidelberg

PHYSIKALISCHES

ΤU

INSTI





Trap modulation spectroscopy

#### **Measuring energies**

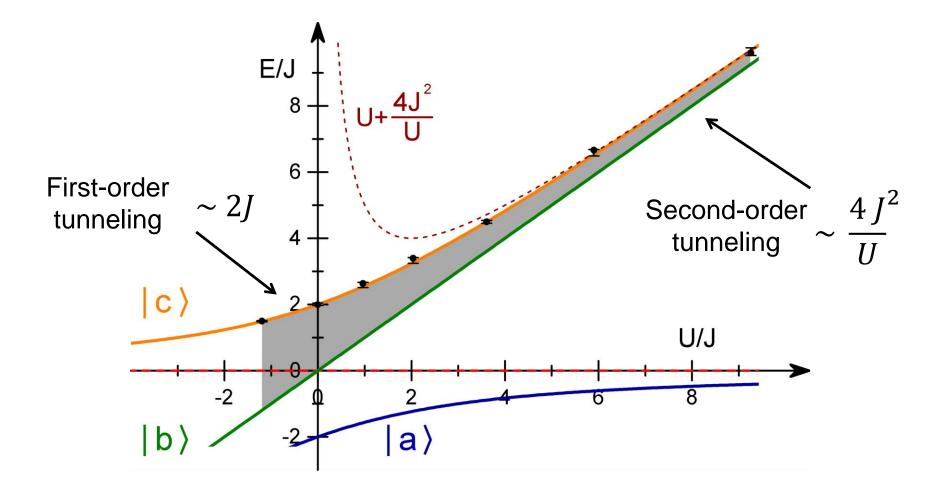
Universität Heidelberg

PHYSIKALISCHES

ΤU

INSTI



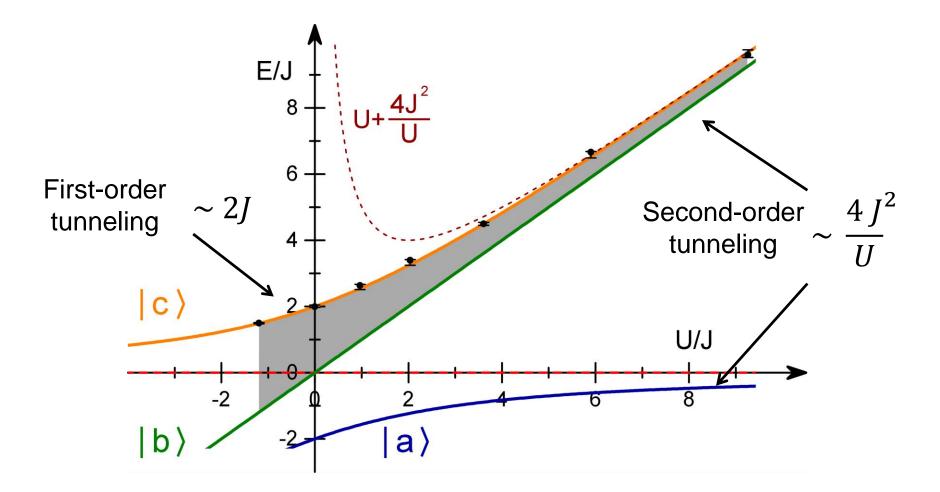


Trap modulation spectroscopy

### **Measuring energies**

Universität Heidelberg





Super exchange energy! responsible for spin ordering in the many body ground state

PHYSIKALISCHES INSTITUT



Universität

Heidelberg

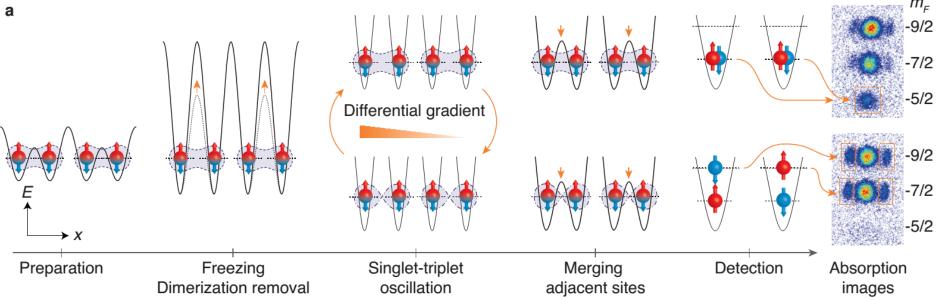
PHYSIKALISCHES

ΙN

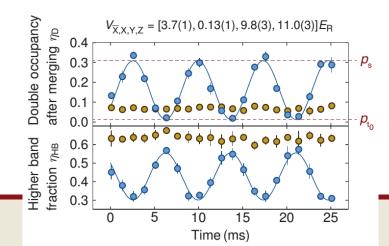
S

# How to go to a many-body system?

• Inspired by a top-down approach: D. Greif et al., Science 340, 1307-1310 (2013) (ETH Zürich)



Dimerize a lattice filled with spin-1/2 fermions to observe spin correlations



# Outlook



Combination of multiple double wells

 Preparation of ground states in separated double wells

Combination to larger system

Can this process be done adiabatically ? Can it be extended to larger systems ?







Dhruv

Kedar



Mathias Neidig

Sebastian Pres

**Puneet Murthy** 

Andrea Bergschneider

Selim Jochim Gerhard Zürn Vincent Klinkhamer

Simon Murmann

# Thank you for your attention!

**Martin Ries** 

Funding:

Center for Quantum Dynamics



