

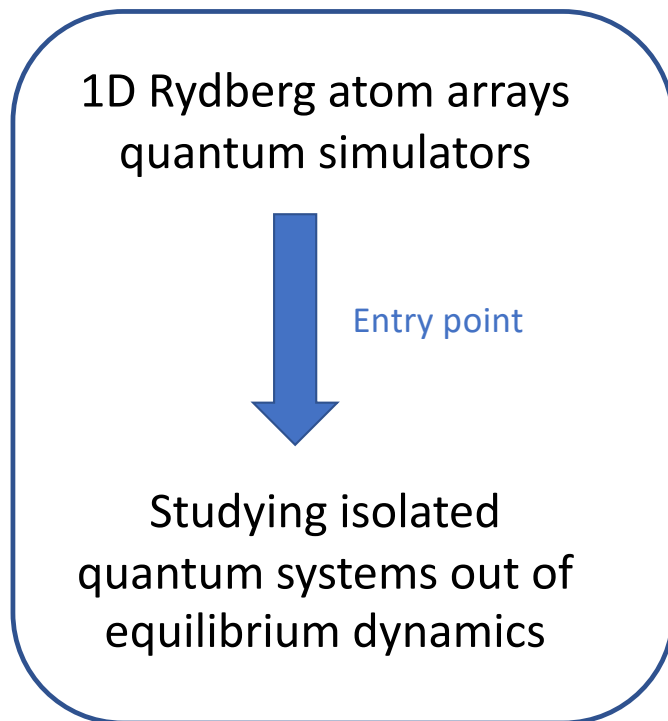


1D Rydberg atom arrays: An entry point to studying isolated quantum many body systems out of equilibrium

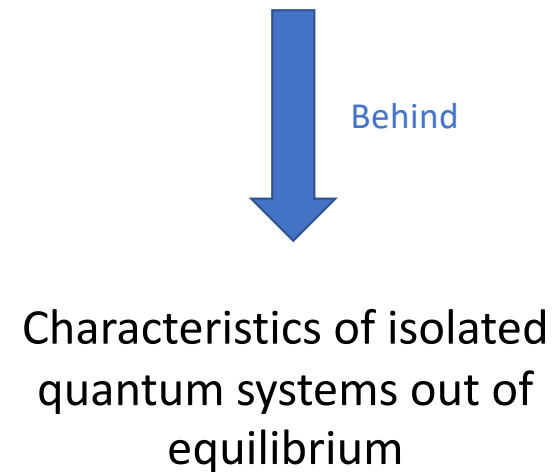
Name: Oliver Lind
Supervisor: Prof Stuart Adams

Quick Aside

Two ideas for this talk

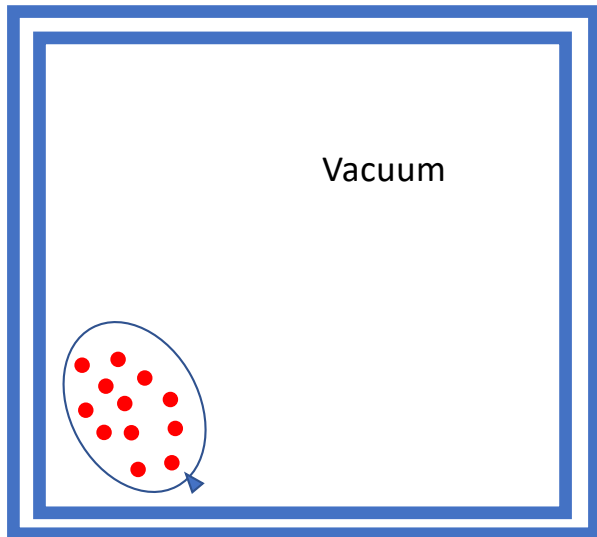


Entanglement entropy dynamics
of system



Motivation

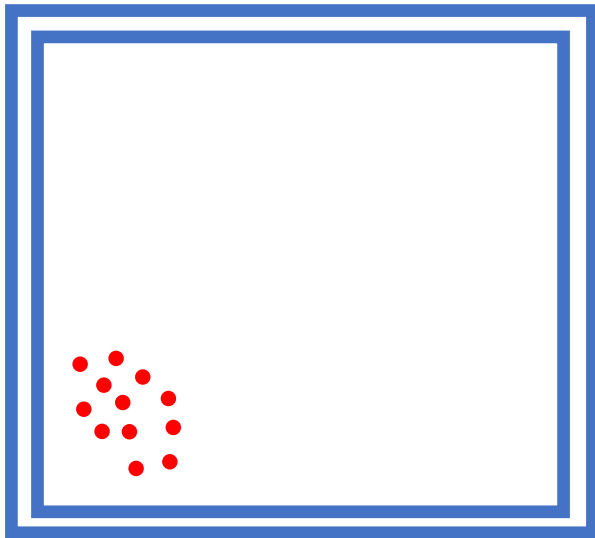
Classical Isolated System



- Adiabatic walls (no transfer of heat outwards)
- Balloon with ideal gas inside
- Rest of the chamber is a vacuum

Motivation

Classical Isolated System

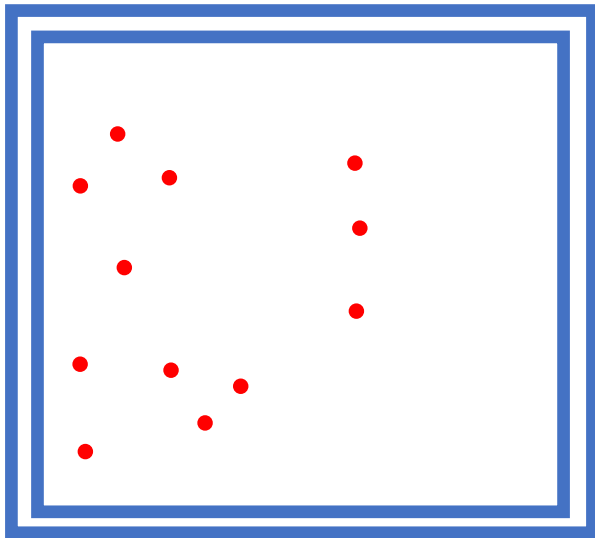


Out of equilibrium initial state

- Adiabatic walls (no transfer of heat outwards)
- Balloon with ideal gas inside
- Rest of the chamber is a vacuum
- **Pierce the balloon**

Motivation

Classical Isolated System

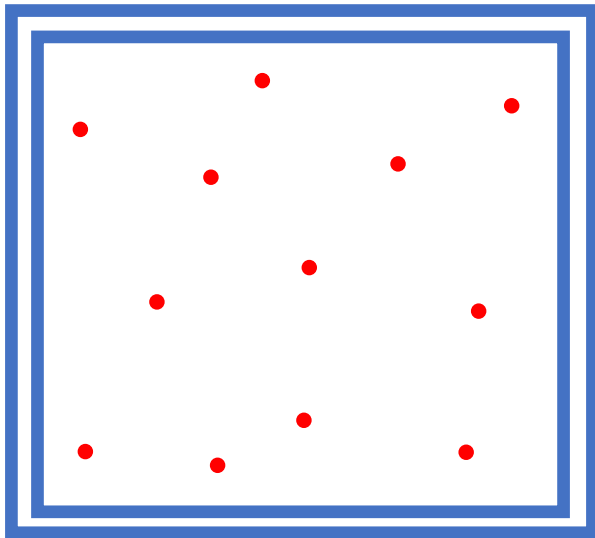


Out of equilibrium state

- Adiabatic walls (no transfer of heat outwards)
- Balloon with ideal gas inside
- Rest of the chamber is a vacuum
- **Pierce the balloon**
- **Air particles spread**

Motivation

Classical Isolated System



Reaches Equilibrium

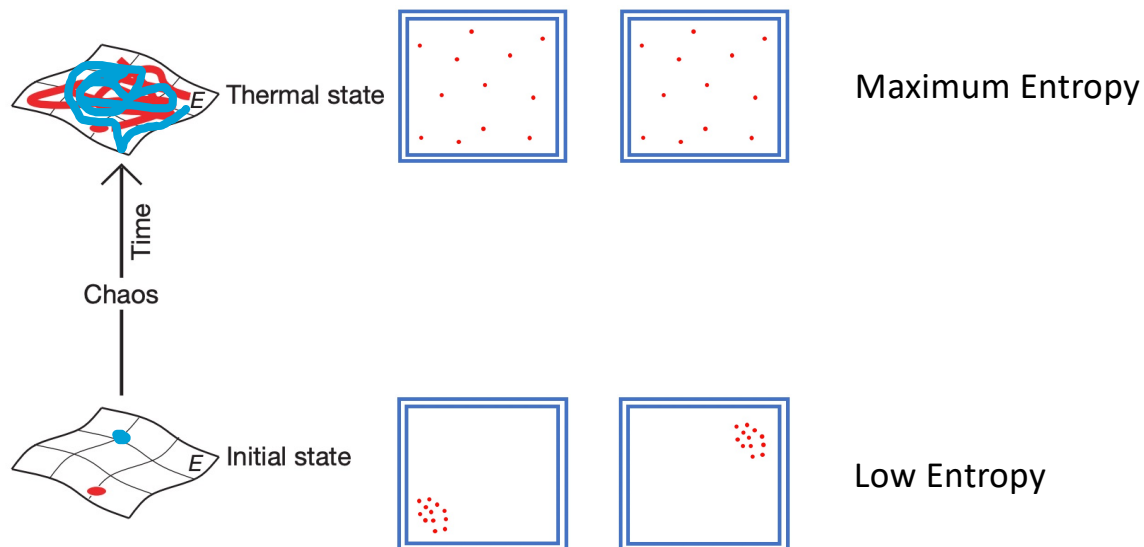
- Adiabatic walls (no transfer of heat outwards)
- Balloon has N particles with total energy E
- Rest of the chamber is a vacuum
- **Peirce the balloon**
- **Air particles spread**
- **System thermalizes**

Macroscopic observables reach steady average value:

$$\langle O(t) \rangle \longrightarrow \langle O \rangle_{thermal}$$

Motivation

Statistical Physics



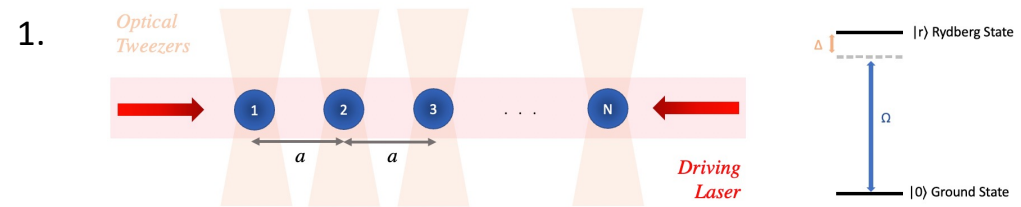
What happens in an isolated quantum system?

- Unitary time evolution
$$|\varphi(t)\rangle = U(t)|\varphi(0)\rangle$$

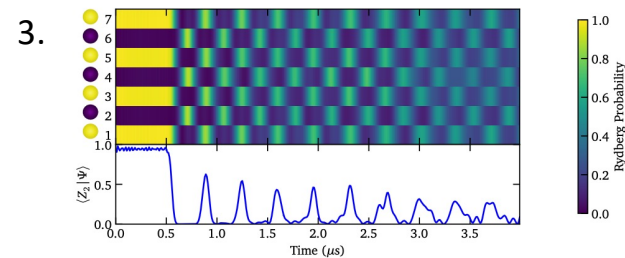
- Can the system thermalize in the same way?
- Can we associate entropy in the same way?
- Are there global properties that can help us understand this?

Contents

1. Introduce the model
 - Rydberg Atoms
 - Rydberg Blockade
 - Ising Type Hamiltonian

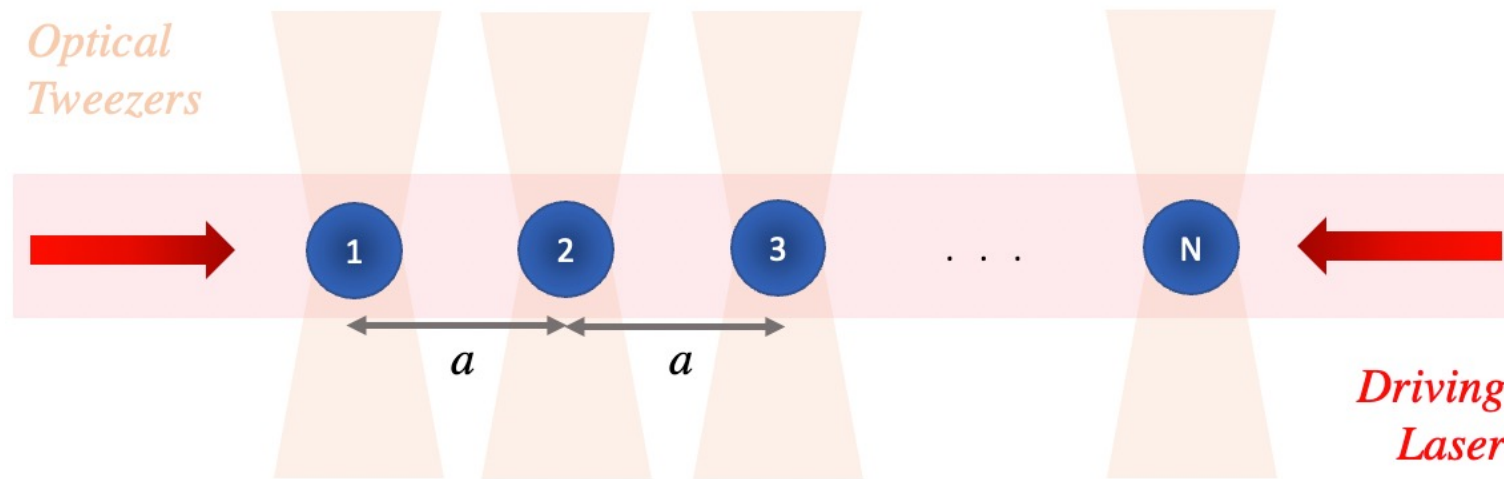


2. Generating order phases
3. Bringing the system out of equilibrium
4. Why is it a good entry point?



Running numerical simulations of quantum simulation protocols done on ultra cold neutral atoms

Set up



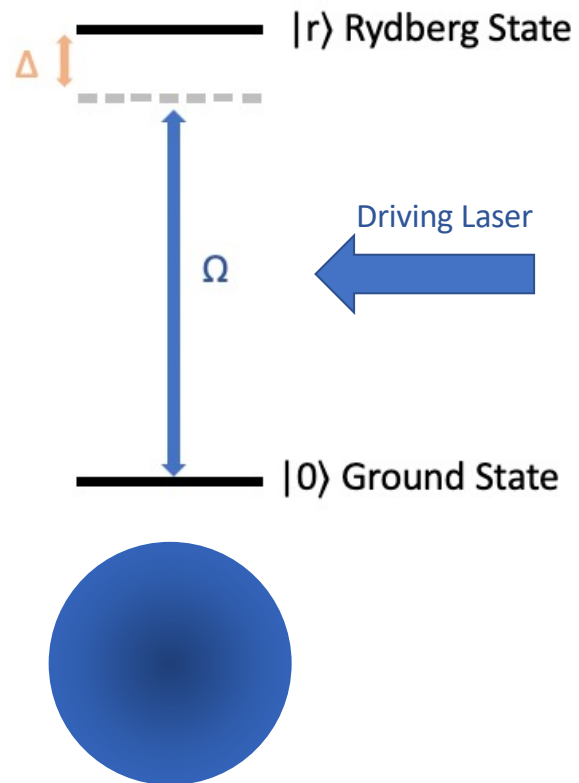
$$\frac{H}{\hbar} = \underbrace{\frac{\Omega(t)}{2} \sum_i^N \sigma_i^x}_{\text{Driving term}} - \underbrace{\Delta(t) \sum_i^N n_i}_{\text{Detuning term}} + \underbrace{\sum_{i < k}^N V_{ik} n_i n_k}_{\text{Interaction term}} \quad i = \text{atom site}$$

Rydberg Atoms: What are they?

- Single electron in a highly excited state ($n \sim 40 - 100$)
- Appears Hydrogen-like
- 'Giant Atoms'
- Large dipole moment
- Strong van der Waals interactions
- Highly excited state is called the Rydberg state



Rydberg Atoms: Encoding as qubit

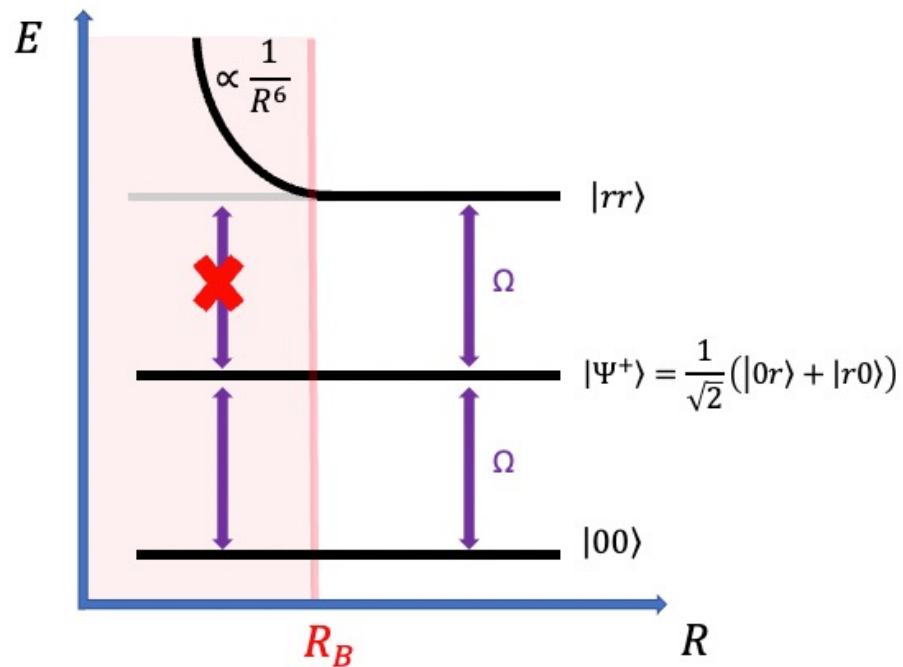


- Use a laser to couple Ground and Rydberg states of atom
- Two level system forms our qubit

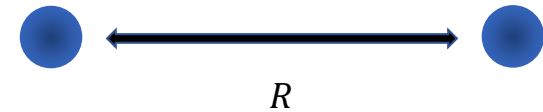
$\frac{\Omega}{2\pi}$ - Rabi Frequency (MHz) Strength of coupling between states

$\frac{\Delta}{2\pi}$ - Detuning (MHz) How far of resonance driving laser is

Interaction and Rydberg Blockade



Two atoms



- Strong **van der Waals interaction** between Rydberg atoms $\propto 1/R^6$
- This results in energy shift of $|rr\rangle$ state
- The blockade radius R_B – characteristic radius where this starts to take effect
- Shared excitation amongst atoms result in entangled states

Many Body Hamiltonian

Evolve the system with Trotter-Suzuki decomposition:

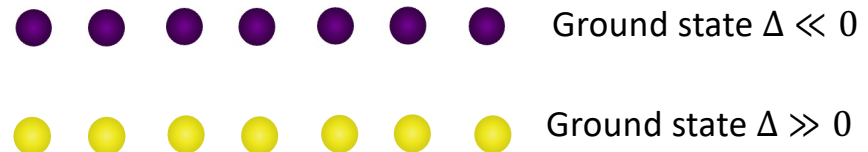
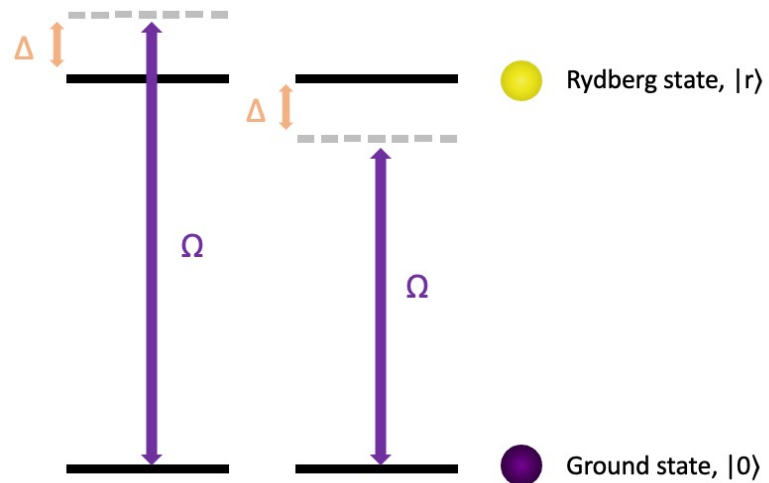
$$|\Psi(t)\rangle = \prod_{i=1}^N e^{-iH[\Omega, \Delta(t_i)]dt_i} |\Psi(t=0)\rangle$$

$$dt_i = 0.01 \mu s$$

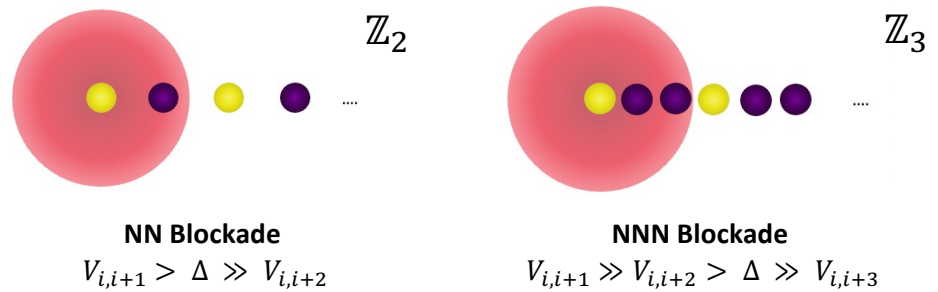
$$\frac{H}{\hbar} = \underbrace{\frac{\Omega(t)}{2} \sum_i^N \sigma_i^x}_{\text{Driving term}} - \underbrace{\Delta(t) \sum_i^N n_i}_{\text{Detuning term}} + \underbrace{\sum_{i < k}^N V_{ik} n_i n_k}_{\text{Interaction term}} \quad i = \text{atom site}$$

$$\Omega(t)/2\pi = \Omega_0/2\pi = 4.00 \text{ (MHz)}$$

Using blockade to generate ordered states



But Rydberg Blockade constraint.... So for $\Delta \gg 0$



Note:

- Focus NN Blockade
- Odd number of sites (No degeneracy in our $\Delta \gg 0$ ground states)

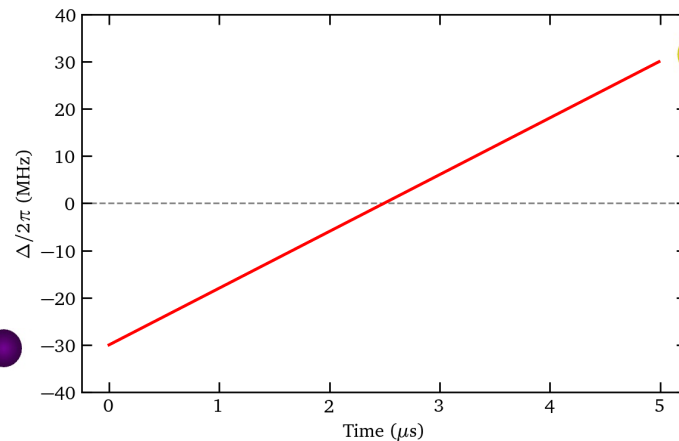
Using blockade to generate ordered states

Evolve the system adiabatically

By adjusting the separation between atoms a , we can sweep through detuning values to generate desired ordered states:

Focus on 7
Atoms

$\Delta \ll 0$ Ground state



$\Delta \gg 0$ Ground State
 \mathbb{Z}_2

Note

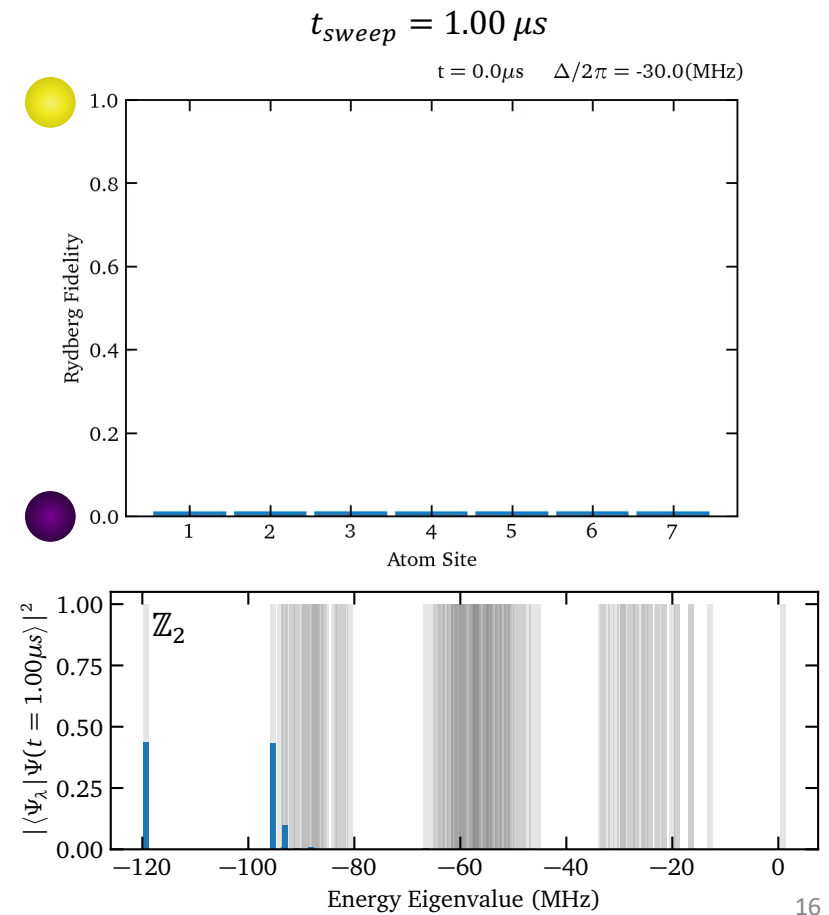
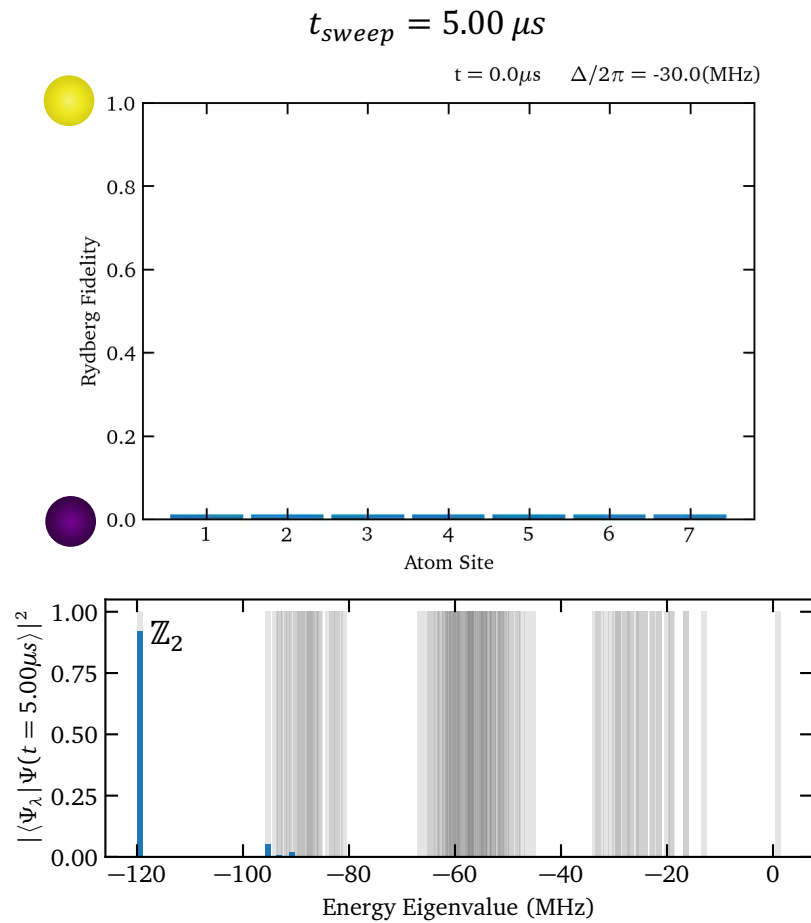
- For
- On

$\Delta \gg 0$

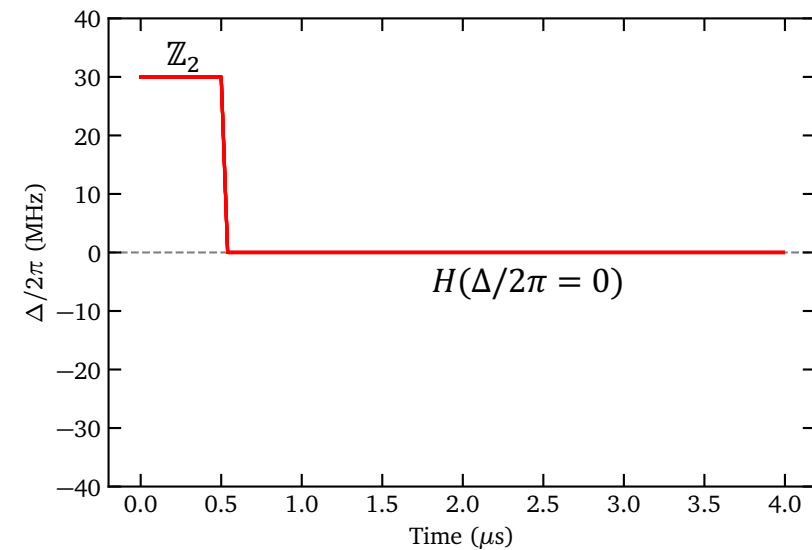
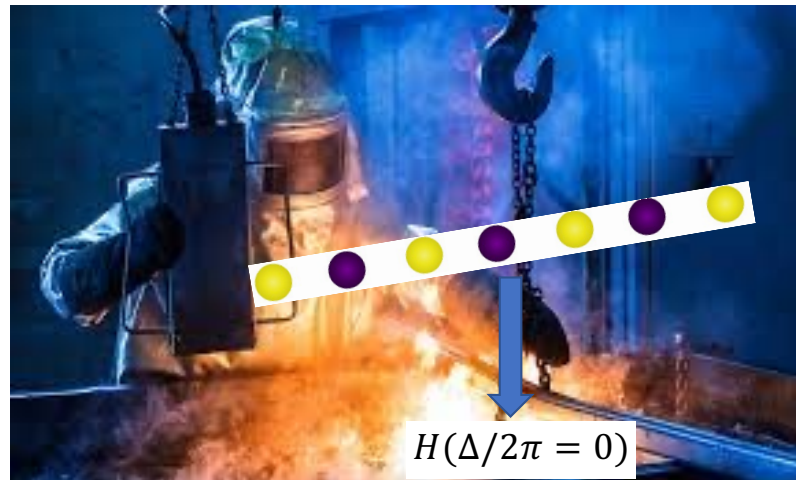
Odd GS

Even GS

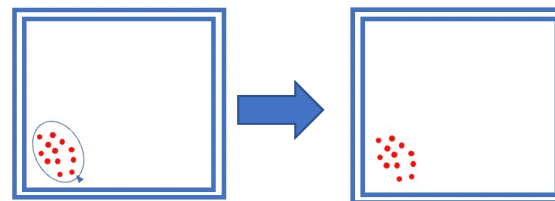
Using blockade to generate ordered states



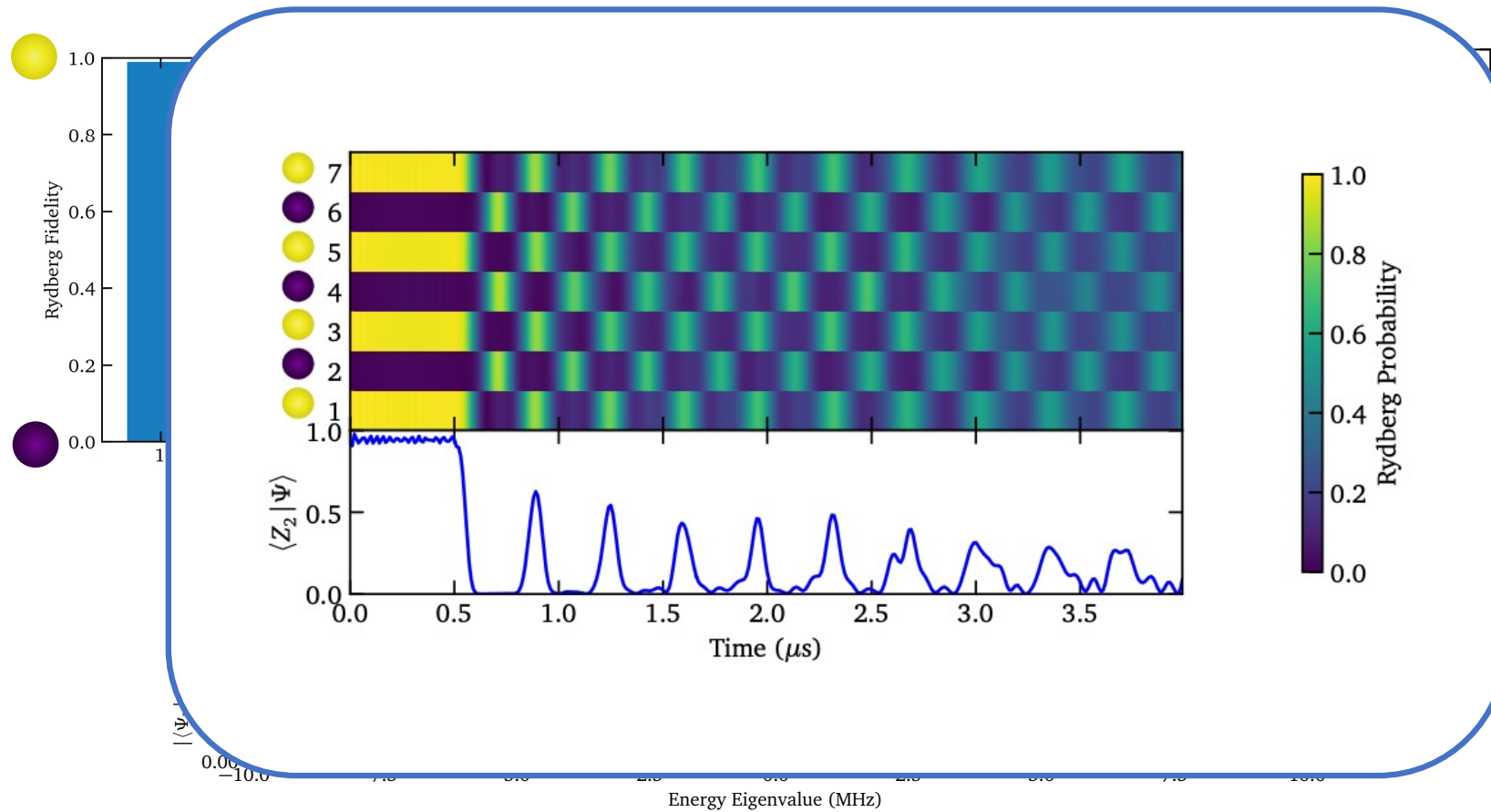
What if we intend on going quickly: quenches



- How we take the system out of equilibrium
- Analogues to balloon popping



Taking the system out of equilibrium

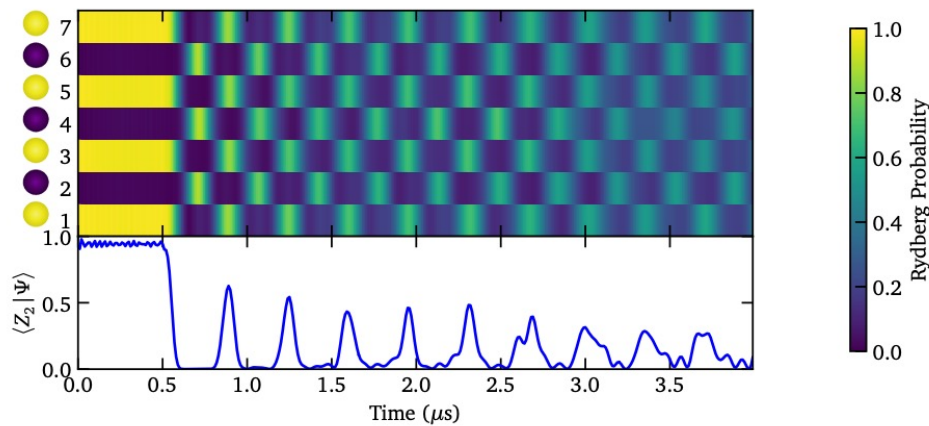


So why is this a good entry point?

1. Interesting novel behavior
 - Quantum many body scars
2. Tunability and versatility to try different things
 - Local quenches
 - Slower quenches
 - Different initial configurations
3. What lies underneath
 - Entanglement
 - Propagation of information

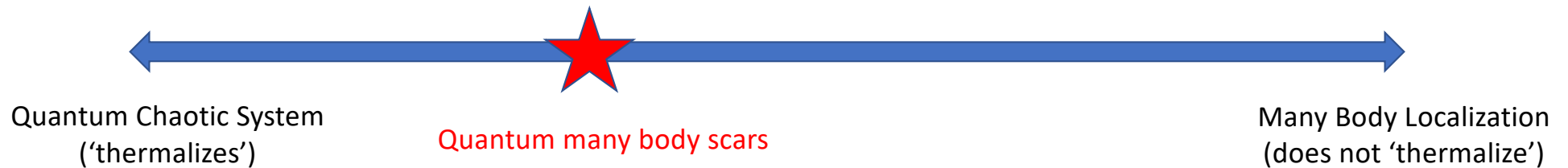
So why is this a good entry point?

1. Interesting novel behavior



- Revival of the initial state after the quench is a behavior known as **quantum many body scars**
- Type of behavior known as **weak ergodicity breaking**

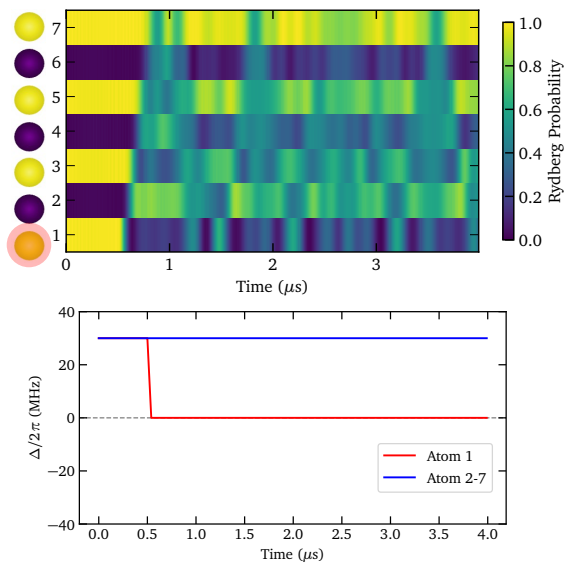
$$|\varphi(t)\rangle = \sum_n a_n e^{-iE_n t} |n\rangle$$



So why is this a good entry point?

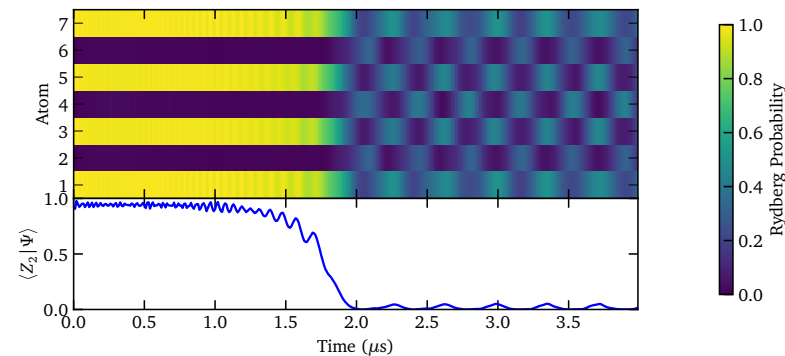
2. Tunability and versatility to try different things

Local addressing and quenches



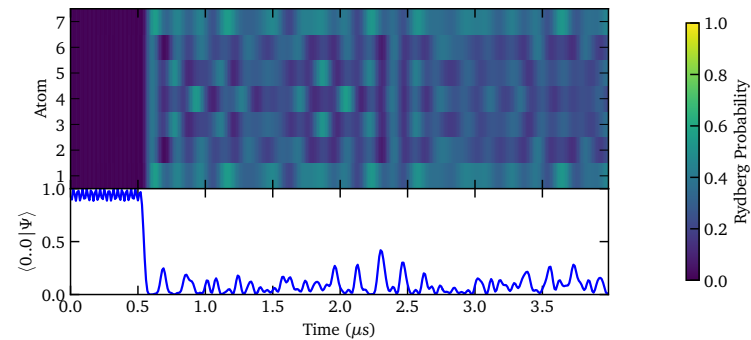
- Propagation of information
- Formations of correlations

Vary the speed to the quench



- Evaluate different out of equilibrium regimes

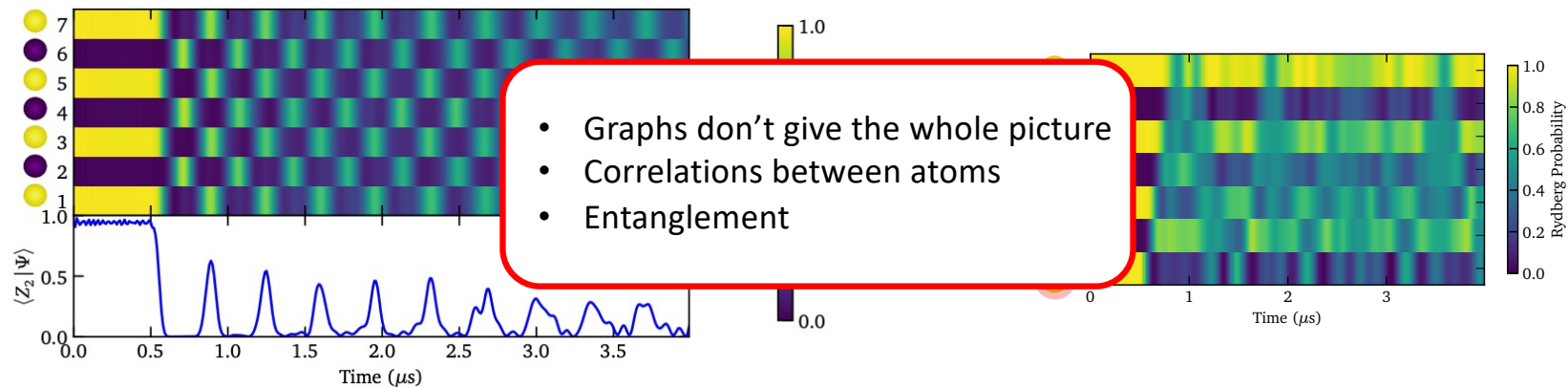
Quench different initial states



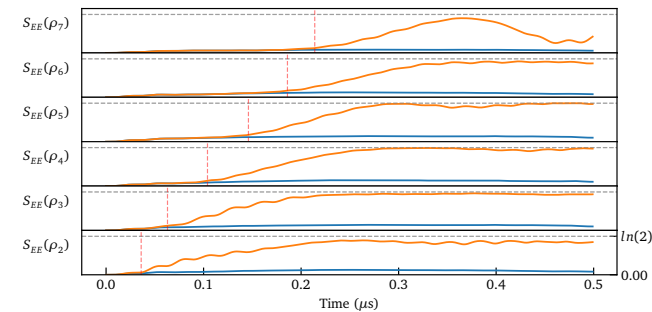
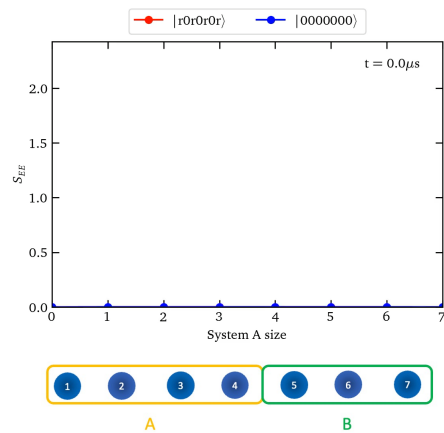
- Investigate which states have scar behavior

So why is this a good entry point?

- What lies underneath



Looking at
entanglement
growth and scaling
with the system



Acknowledgements

- Supervisor: Prof Stuart Adams
- PHD Student: Toonyawat Angkhanawin

Values

$$\frac{H}{\hbar} = \underbrace{\frac{\Omega}{2} \sum_i^N \sigma_i^x}_{\text{Driving term}} - \underbrace{\Delta \sum_i^N n_i}_{\text{Detuning term}} + \underbrace{\sum_{i < k}^N V_{ik} n_i n_k}_{\text{Interaction term}} \quad i = \text{atom site}$$

$$\sigma_i^x = |0_i\rangle\langle r_i| + |r_i\rangle\langle 0_i|$$

$$n_i = |r_i\rangle\langle r_i|$$

$$V_{ik} = \frac{C_6}{r^6}$$

$$\Omega/2\pi = 4.00 \text{ (MHz)}$$

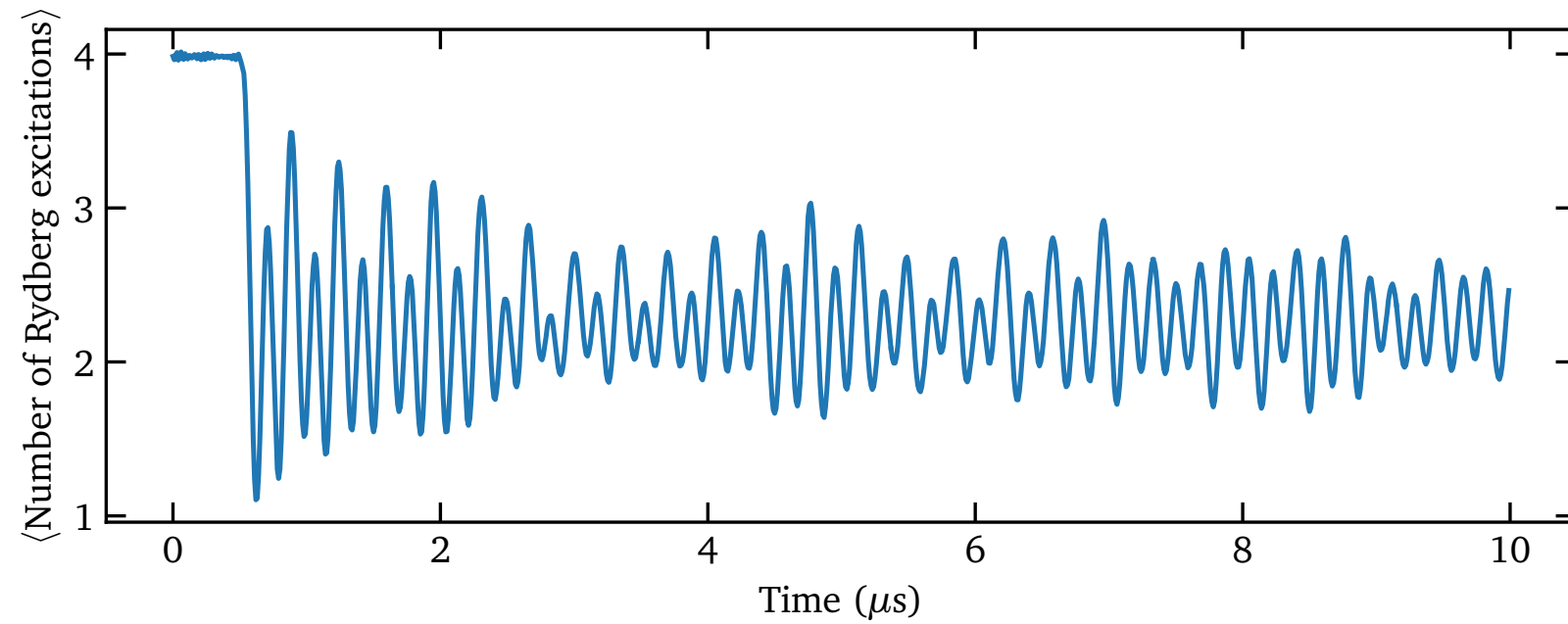
$$r = a|i - k|$$

$$\text{Rb Atoms: } C_6/2\pi = 863000 \text{ MHz } \mu\text{m}^6$$

Evolve the system with Trotter-Suzuki decomposition:

$$|\Psi(t)\rangle = \prod_{i=1}^N e^{-iH[\Omega, \Delta(t_i)]dt_i} |\Psi(t=0)\rangle \quad dt_i = t_i - t_{i-1} \quad dt_i = 0.01 \mu\text{s}$$

Thermalization or lack there of



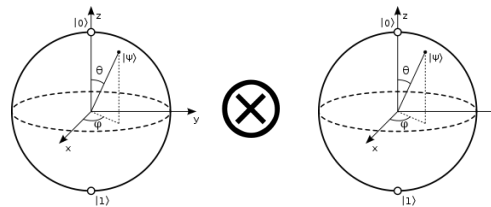
Entanglement: can this offer us any further insight?

Two qubits

Product states

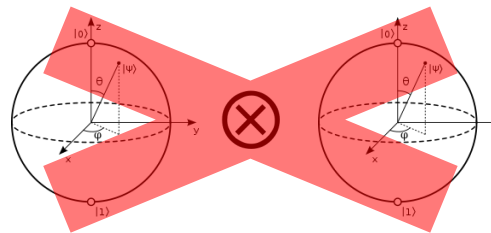
$$\frac{1}{\sqrt{2}}(|0\rangle + |r\rangle) \otimes \frac{1}{\sqrt{2}}(|0\rangle + |r\rangle)$$

Atom 1 Atom 2



Entangled states

$$|\Psi^+\rangle = \frac{1}{\sqrt{2}}(|0r\rangle + |r0\rangle)$$



Need Atom 2 to fully describe the state of atom 1

Entanglement Entropy



$|\varphi\rangle$ Density Matrix $\rho = |\varphi\rangle\langle\varphi|$



$$|\varphi\rangle = |\varphi\rangle_A \otimes |\varphi\rangle_B$$

$$|\varphi\rangle \neq |\varphi\rangle_A \otimes |\varphi\rangle_B$$

Reduced Density Matrix

$$\rho_A = \text{Tr}_B(\rho)$$

Everything we can know about A without B

$$n_i = |r_i\rangle\langle r_i|$$

Von Neuman Entropy Between A and B

How uncertain is the information in ρ_A