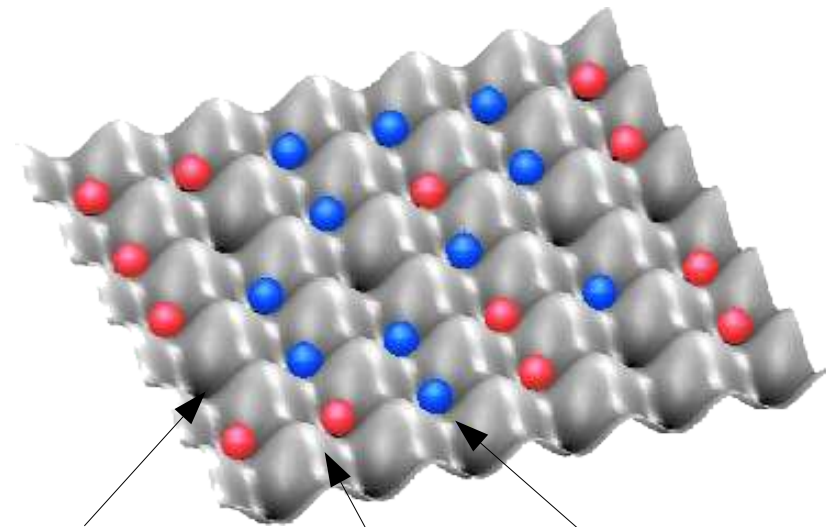


Optical lattice & Quantum computation

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(17-4-2009)

Quantum register



holes

$|0\rangle$

$|1\rangle$

$$N_{1D} \sim 10 - 100$$

$$N_{2D} \sim 100 - 10^4$$

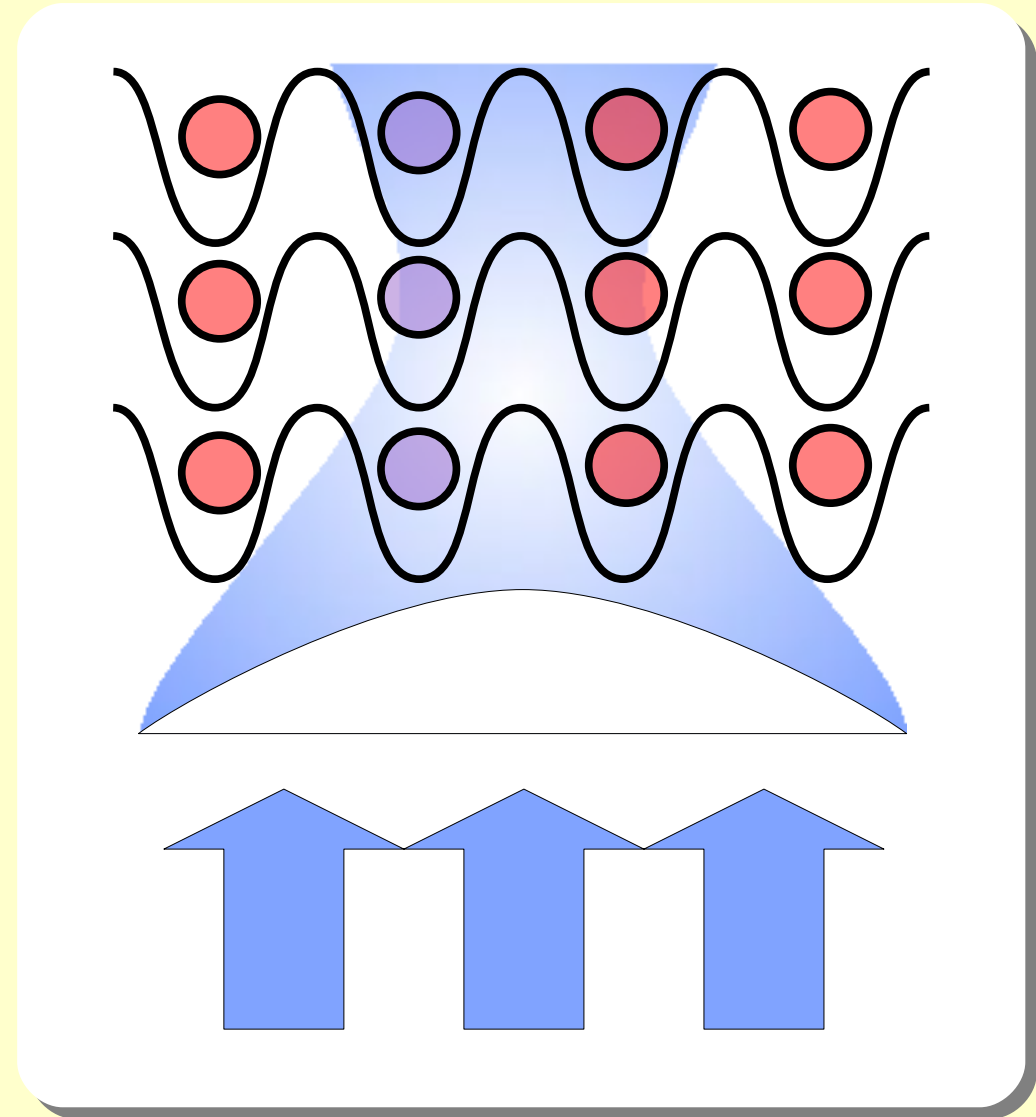
$$N_{3D} \sim 10^6$$

- One atom per lattice site, acting as a possible qubit.
- Information stored in the state of the atom.
- A macroscopic number of qubits...
- ...or a huge number of copies of the same experiment.
- May introduce **defects** when loading the lattice
 - holes, double loading, etc.

Single-qubit operations

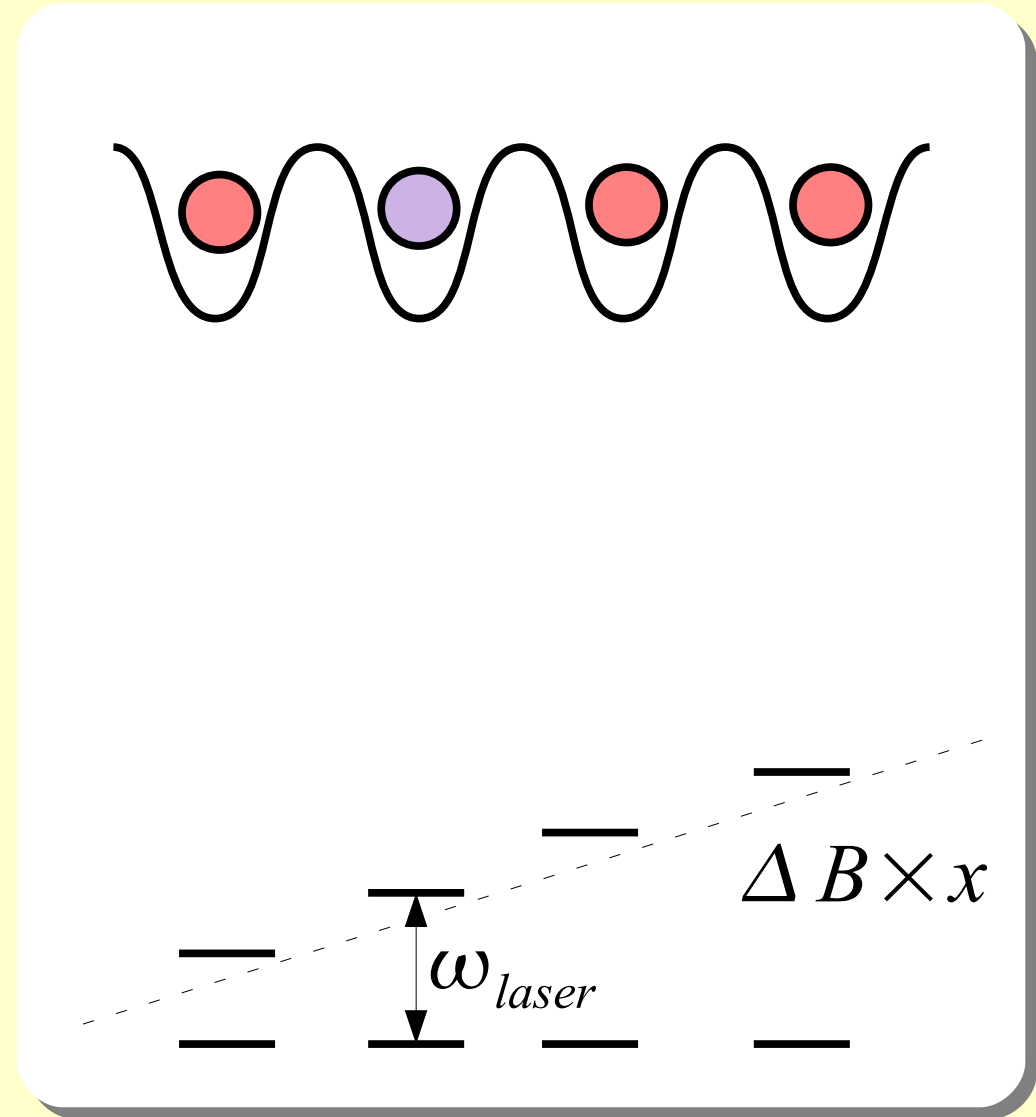
Single-qubit operations

- We have seen how to change the state of qubits.
- But in the lattice it is difficult to achieve individual addressing:
 - manipulate qubits separately
- One reason, diffraction limit:
 - no focusing below wavelength
- Other reason, many layers of atoms in between.



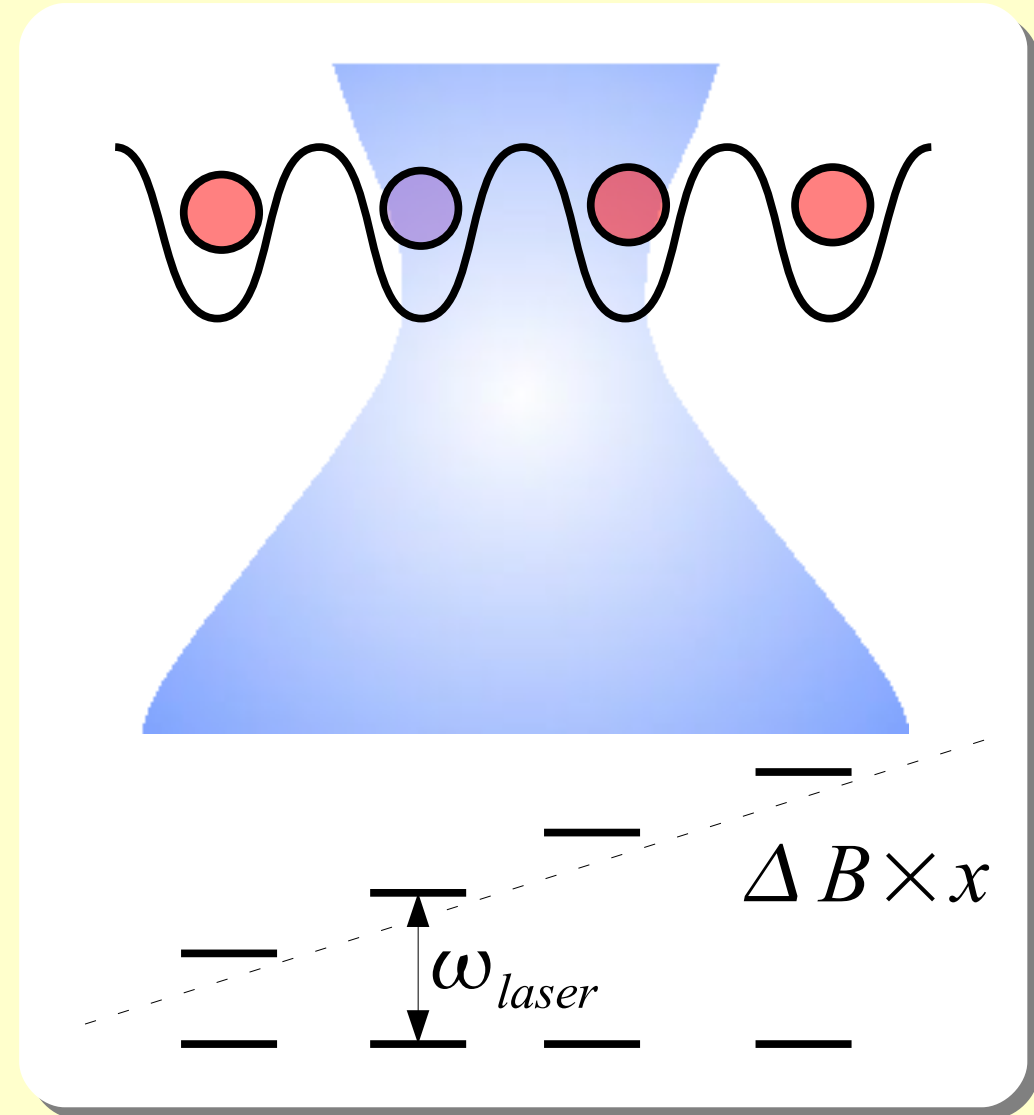
Single-qubit manipulations

- Applying EM field gradients, the energy levels of the atoms change.



Single-qubit manipulations

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- Only some of them become sensitive to the lasers.



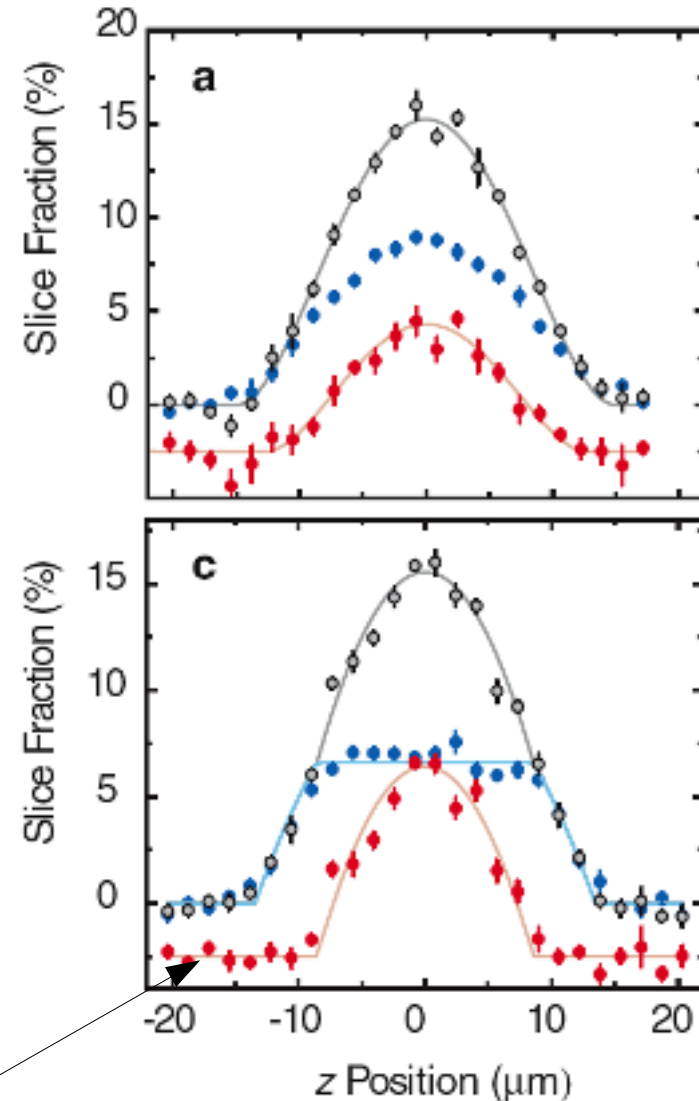
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 - spectrally resolved qubits.



Single-qubit manipulations

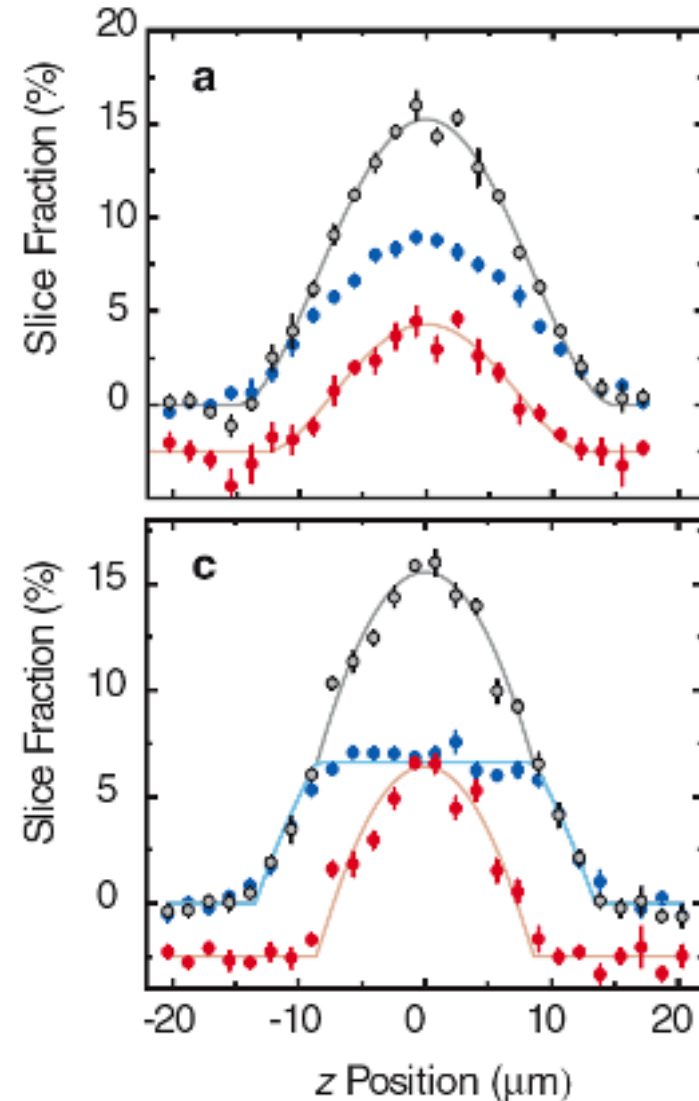
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Each point a single frequency tuning.

Single-qubit manipulations

- Applying EM field gradients, the energy levels of the atoms change.
- Only some of them become sensitive to the lasers.
- Similar in principle to NMR techniques:
 - spectrally resolved qubits.
- But not **enough spatial resolution** yet
 - More ideas later!



Interactions

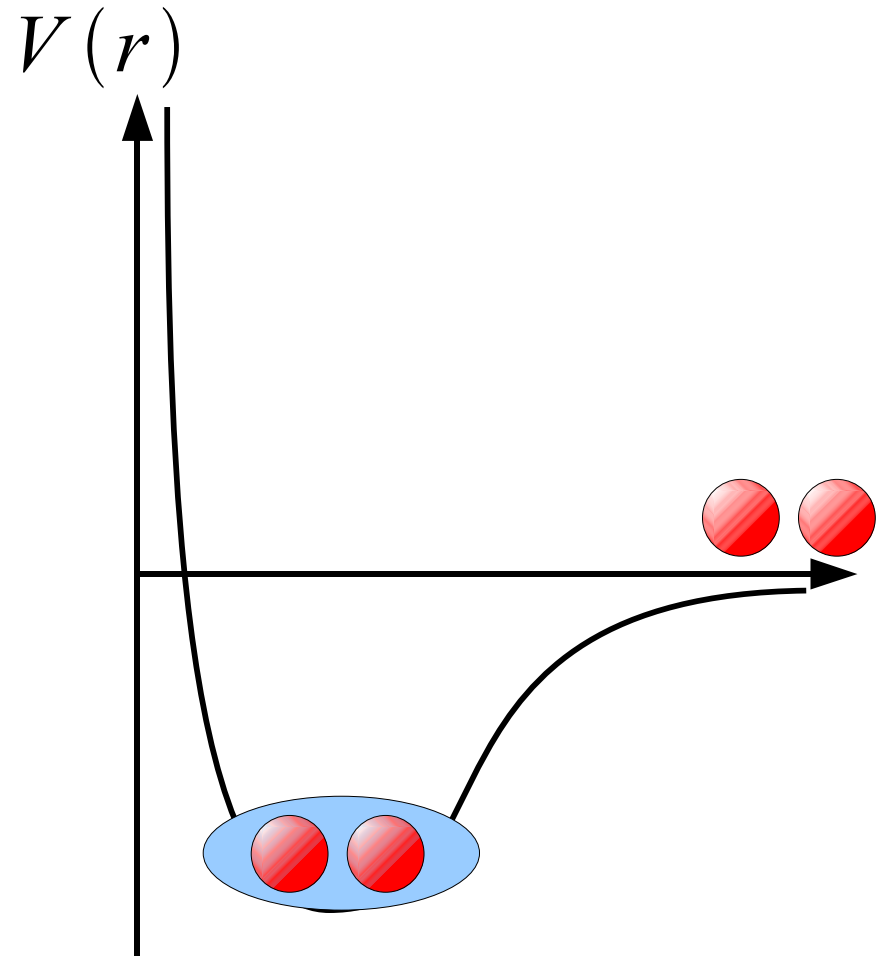
Interactions

- No Coulomb interaction.
- No significant dipolar interaction.
- The state of free atoms is metastable
 - Ground state has molecules or a solid



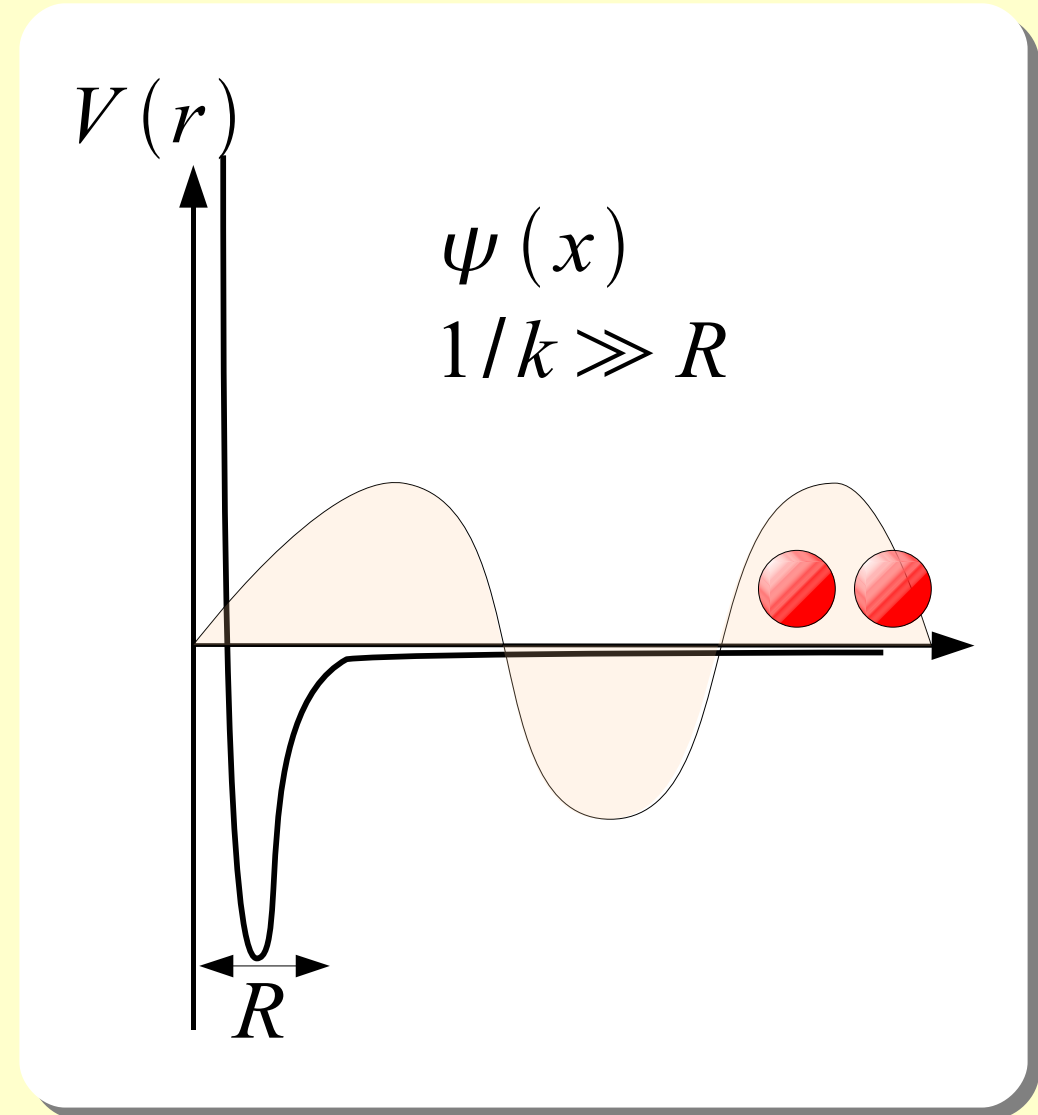
Interactions

- No Coulomb interaction.
- No significant dipolar interaction.
- The state of free atoms is metastable
 - Ground state has molecules or a solid
 - There is some short-range binding potential
 - Too much energy to get actually bound.



Interactions

- When atoms have little kinetic energy, they do not approach much.



Interactions

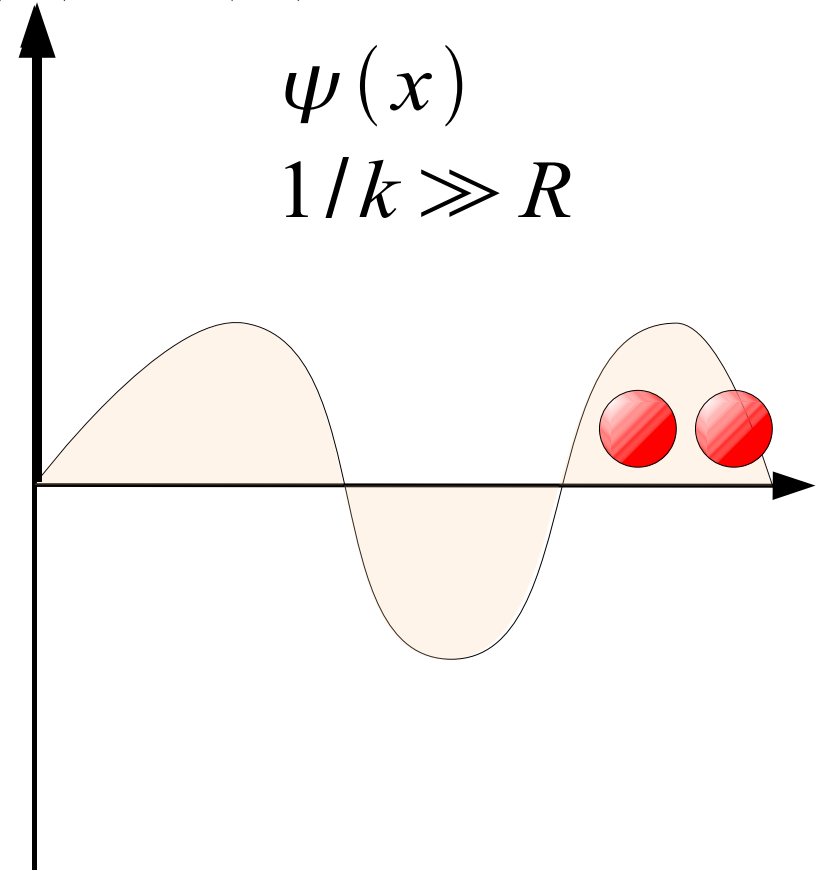
- When atoms have little kinetic energy, they do not approach much.
- We can replace the potential by one with similar scattering properties

$$V(r) \sim \frac{4\pi\hbar^2 a_0}{m} \delta(r)$$

$$V(\vec{r}) \sim \delta(\vec{r})$$

$$\psi(x)$$

$$1/k \gg R$$



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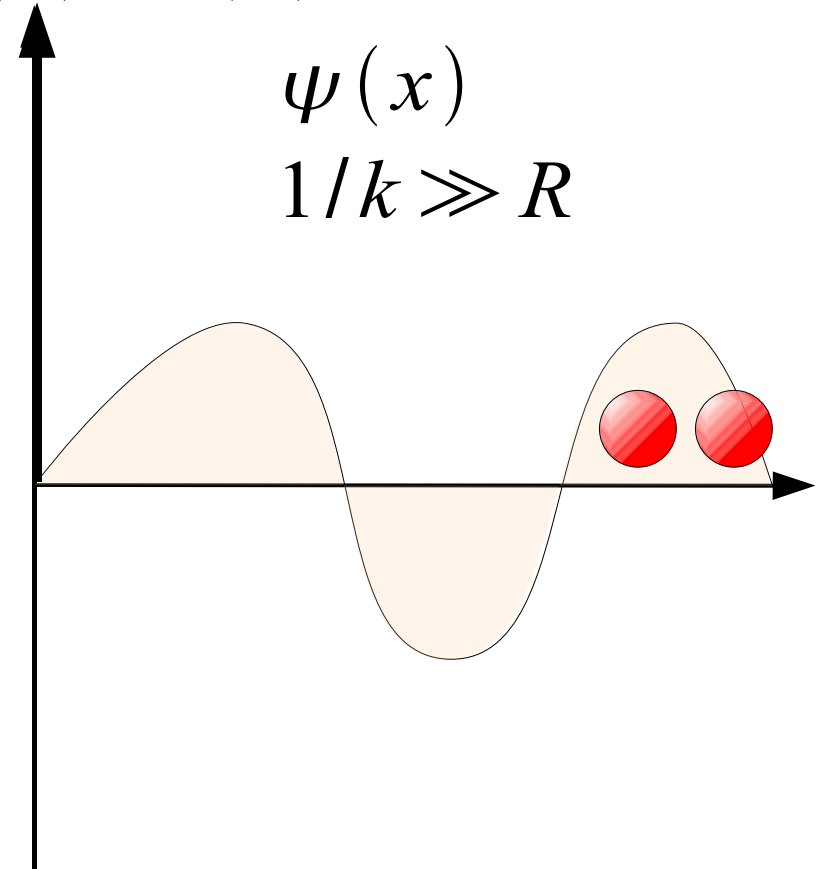
- Second quantization language

$$\frac{g}{2} \int \psi^\dagger(x)^2 \psi(x)^2 dx$$

$$V(\vec{r}) \sim \delta(\vec{r})$$

$$\psi(x)$$

$$1/k \gg R$$



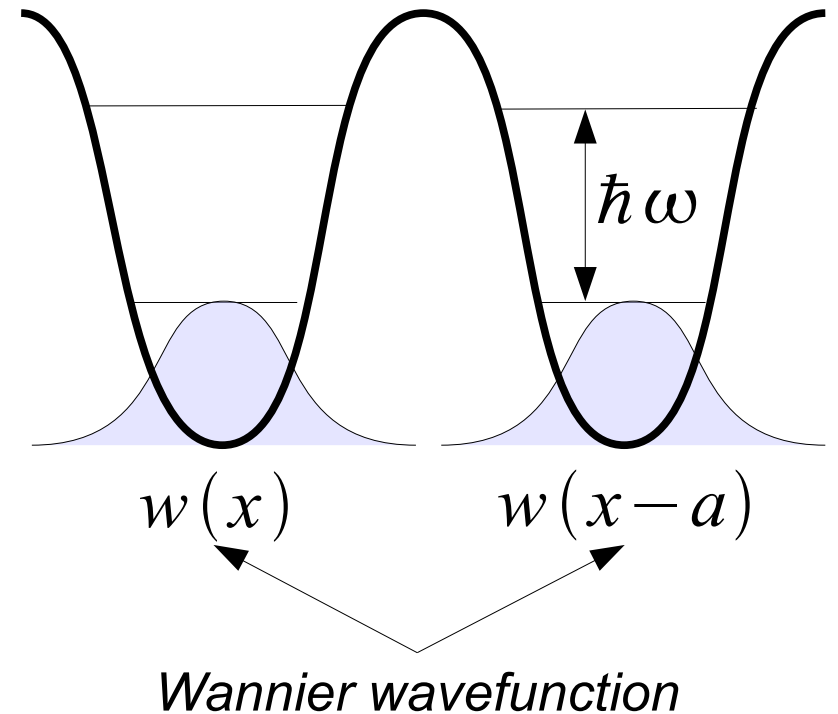
Interactions in the lattice

- Atoms are localized in the approximate ground states of the lattice wells.
- **Tight binding** approximation

$$\psi^+(x) = \sum_i a_i^+ w(x - x_i)$$

- In the interaction we get

$$\begin{aligned} H_{int} &\sim \sum_i \frac{U}{2} a_i^{+2} a_i^2 \\ &= \sum_i \frac{U}{2} n(n-1) \end{aligned}$$

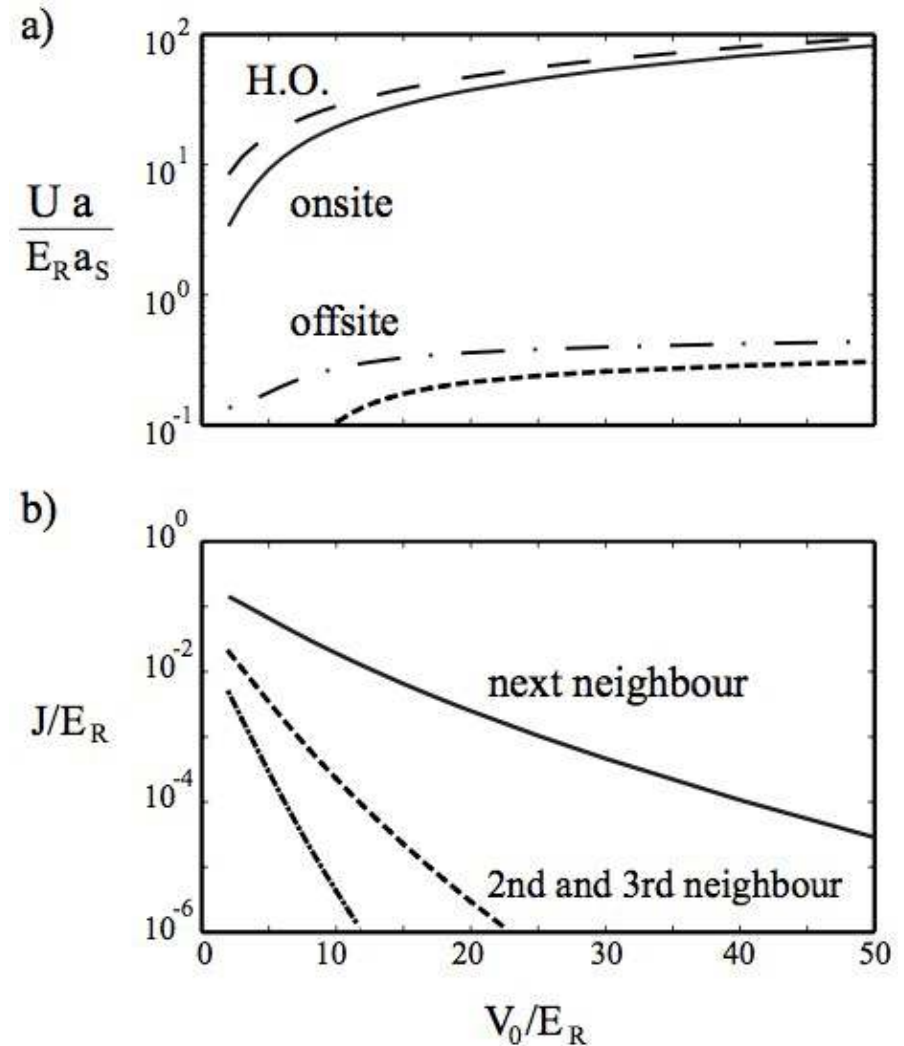


Interactions in the lattice

- The on-site interaction is dominant

$$H_{int} \sim \sum_i \frac{U}{2} a_i^{+2} a_i^2$$
$$= \sum_i \frac{U}{2} n(n-1)$$

- It changes very slowly with the trapping potential.
- Other interaction terms, between neighbors, are negligible.



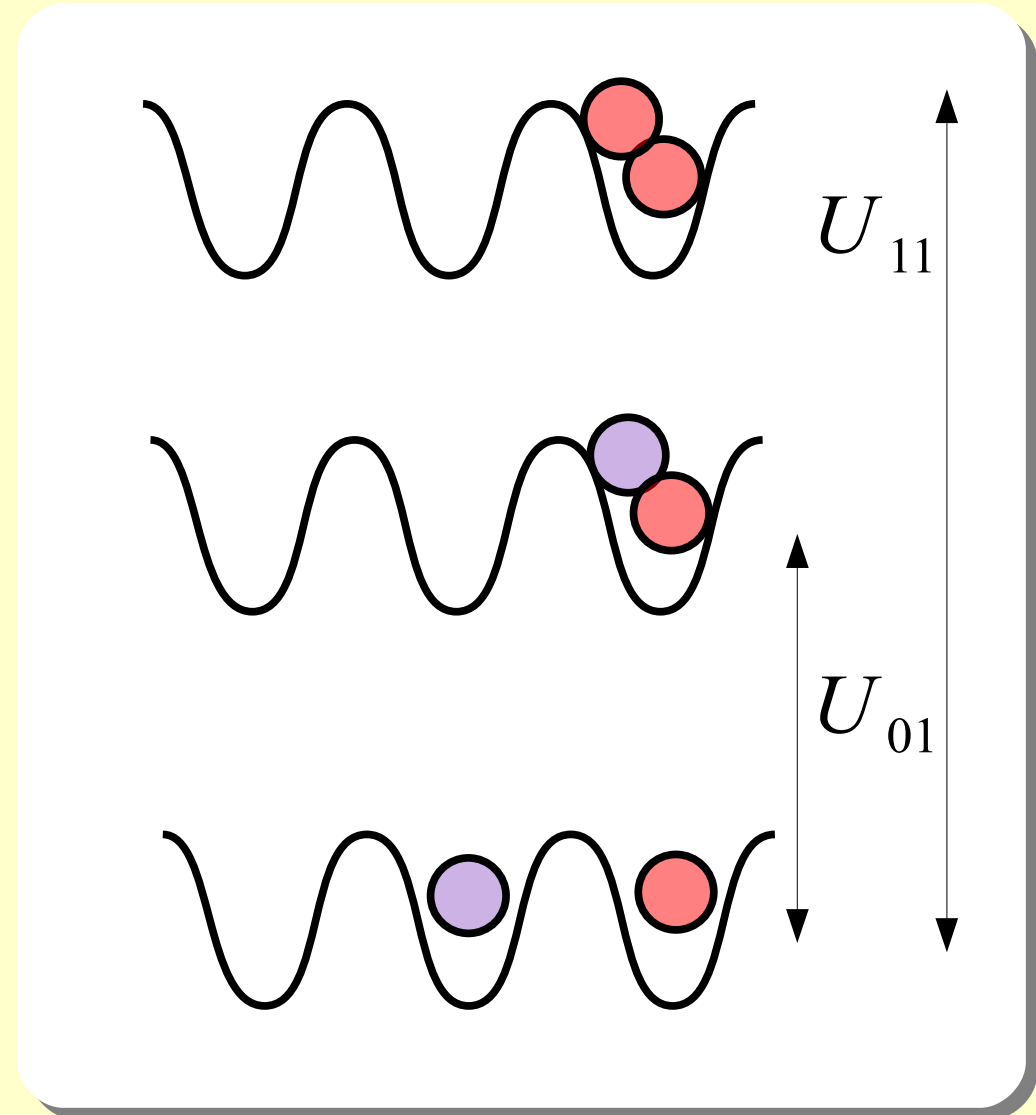
Interactions in the lattice

Summing up:

- Atoms in the same lattice site interact.
- Interaction is “contact”, and too weak to change wavefunction
- Interaction may depend on internal state

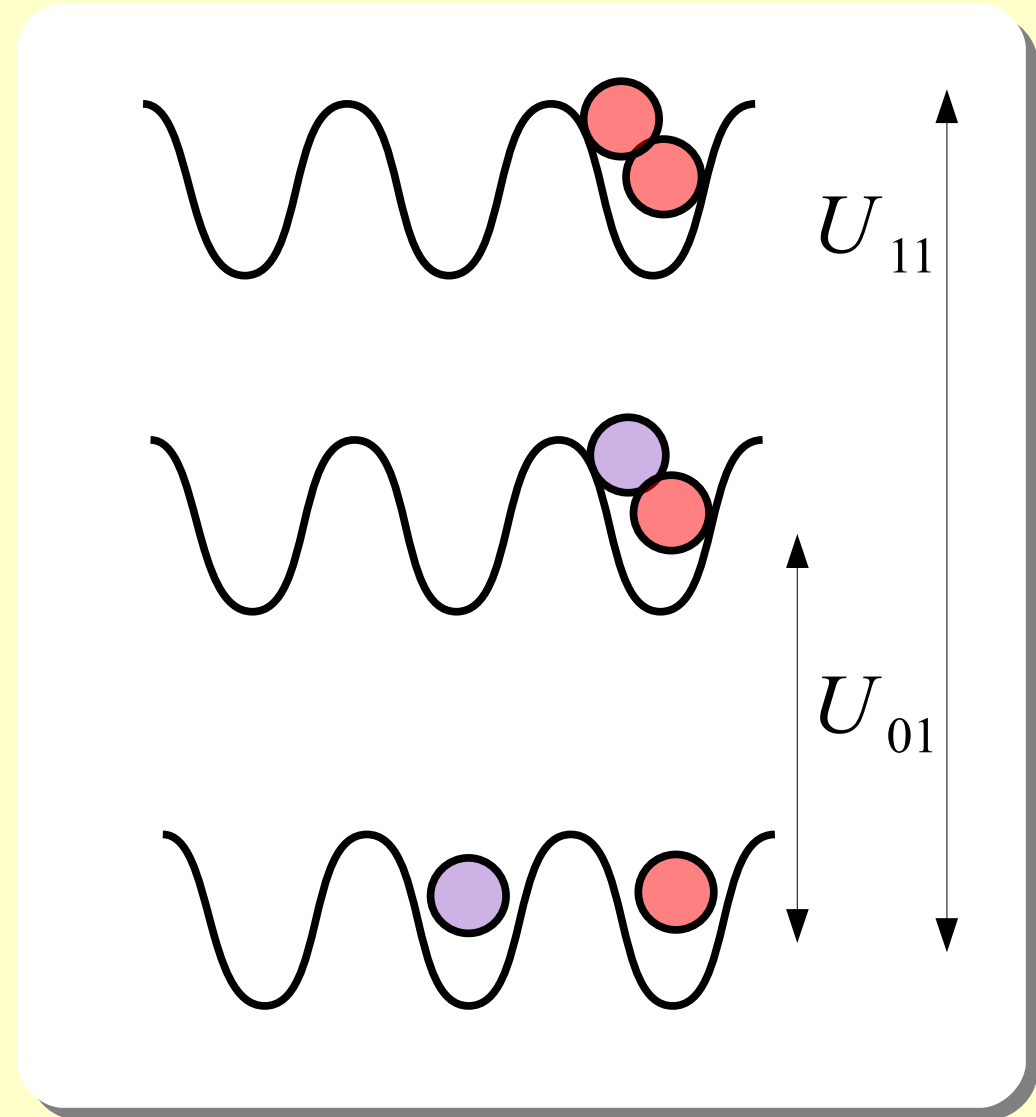
$$H \sim \sum_{\alpha, \beta=0,1} \frac{1}{2} U_{\alpha\beta} n_{\alpha} n_{\beta}$$

- All other terms are negligible



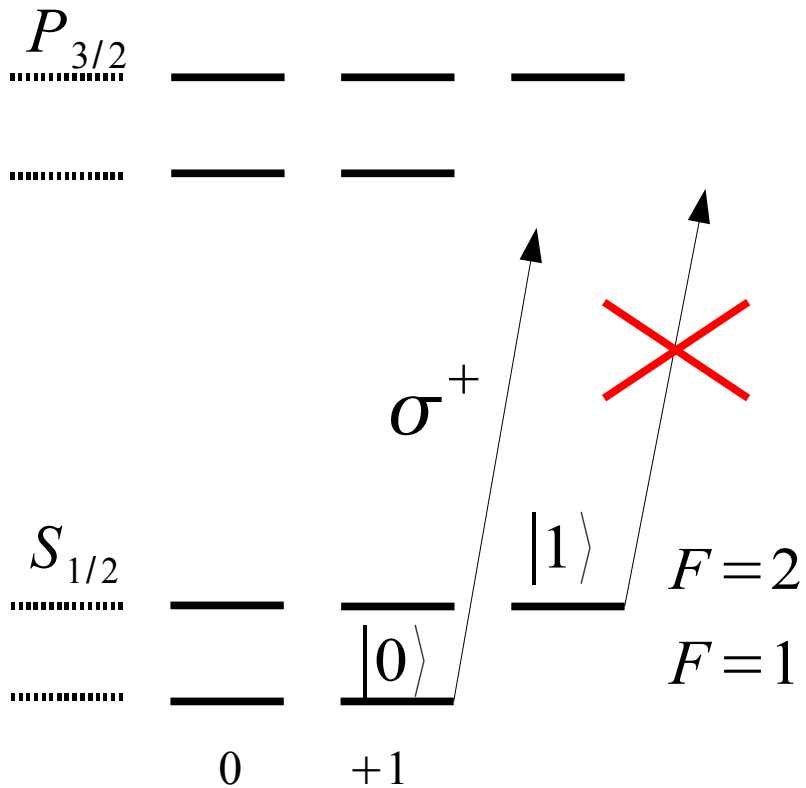
Interactions in the lattice

How do we bring the qubits together to interact?



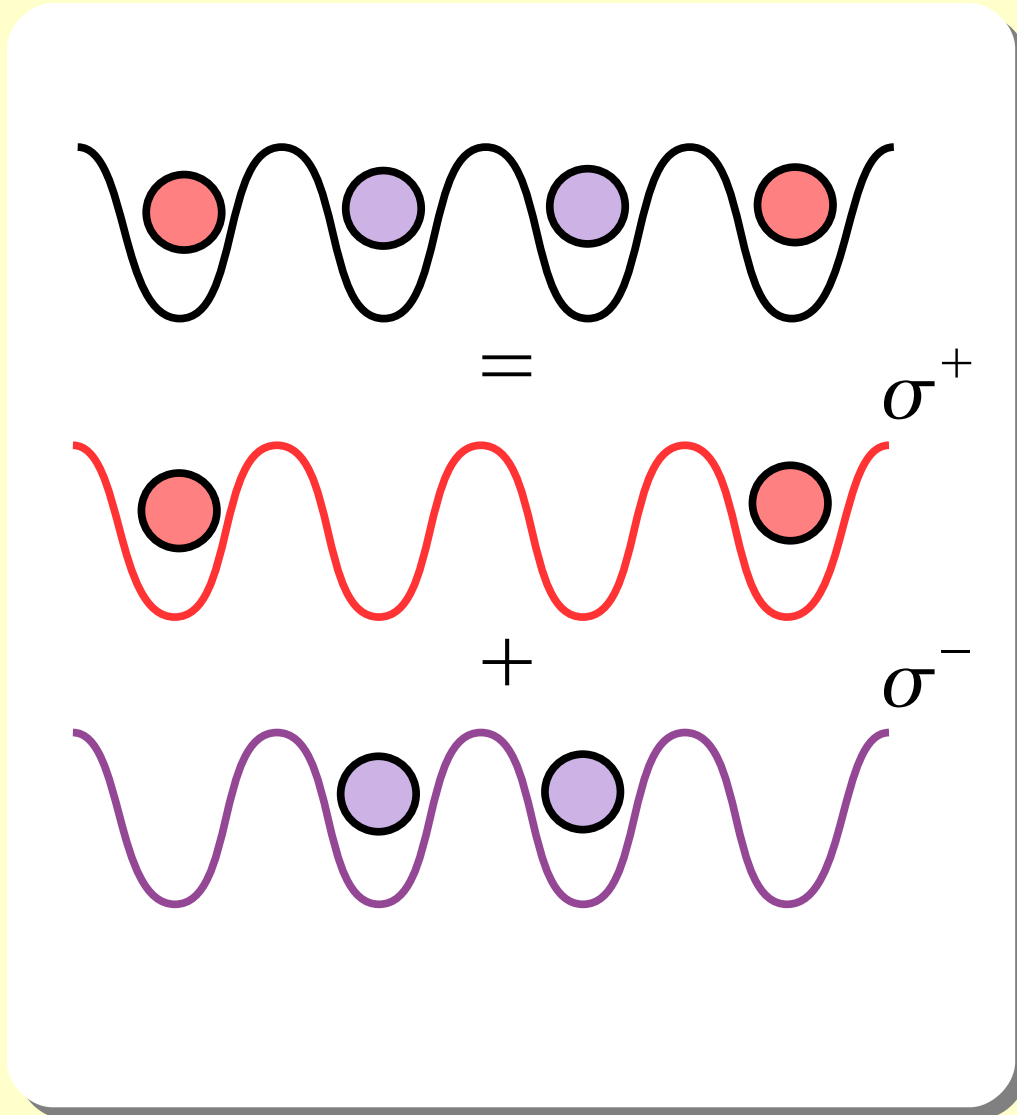
State-dependent lattices

State-dependent lattices



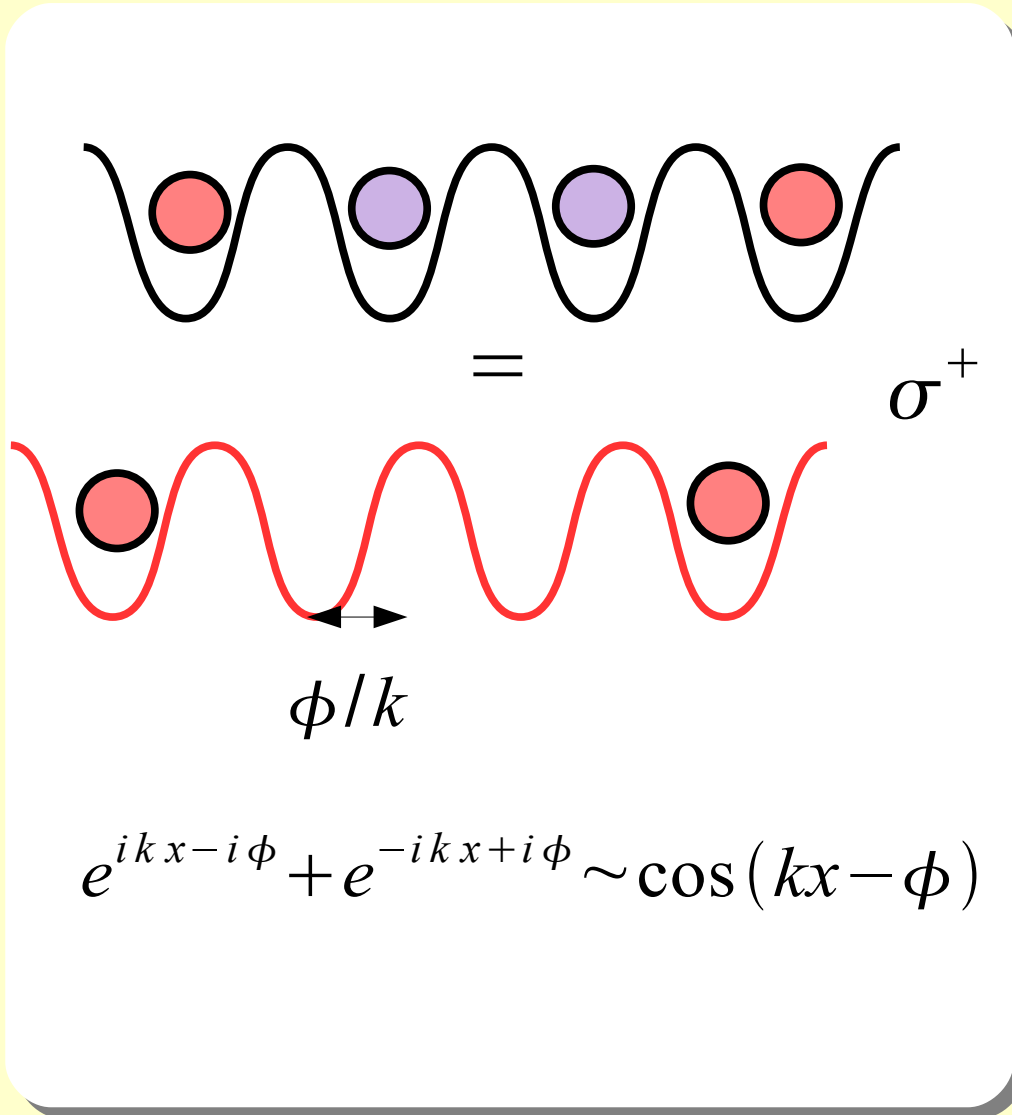
- Not all atomic states react the same to different polarizations.
 - Some transitions are forbidden or suppressed

State-dependent lattices



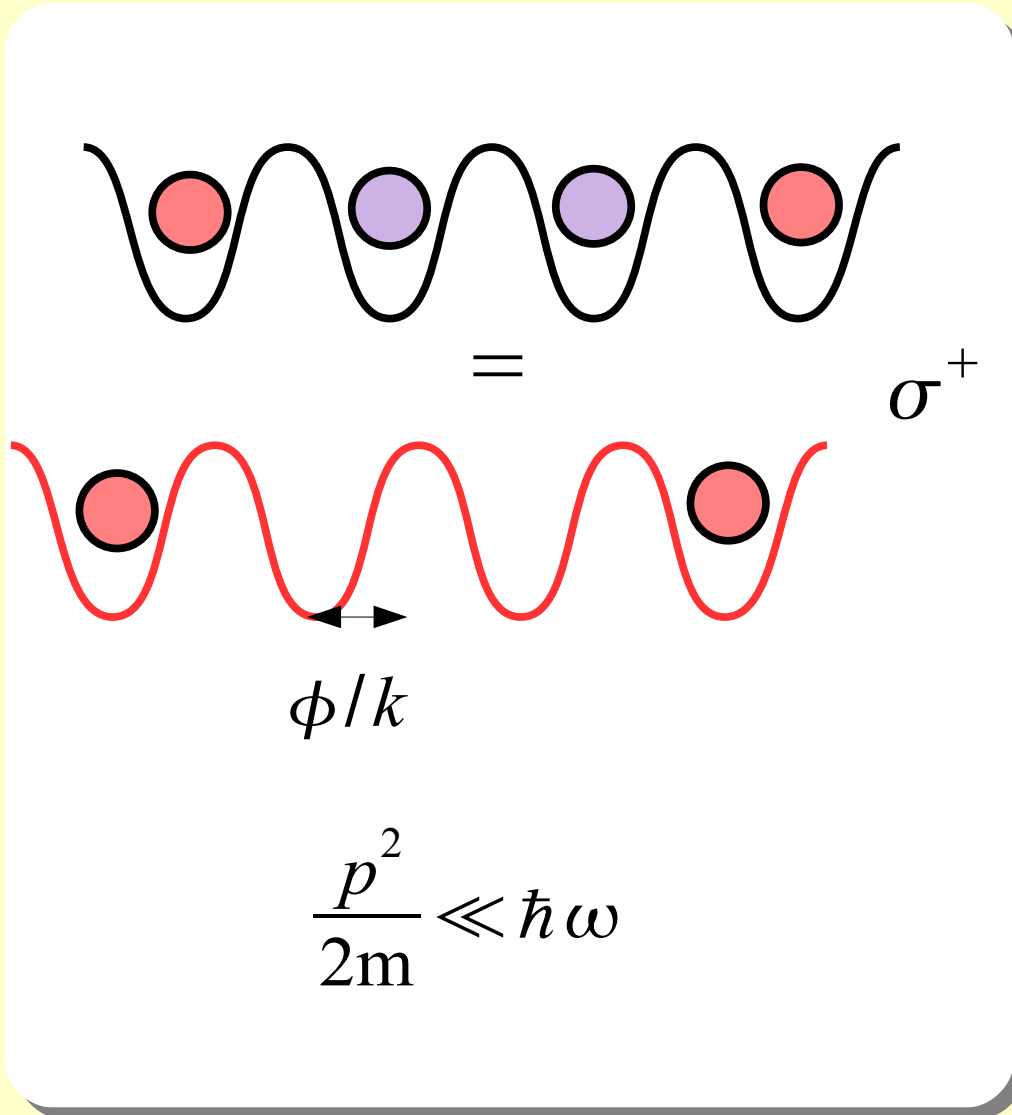
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- We can apply a phase shift (birrefringence) to displace each lattice.

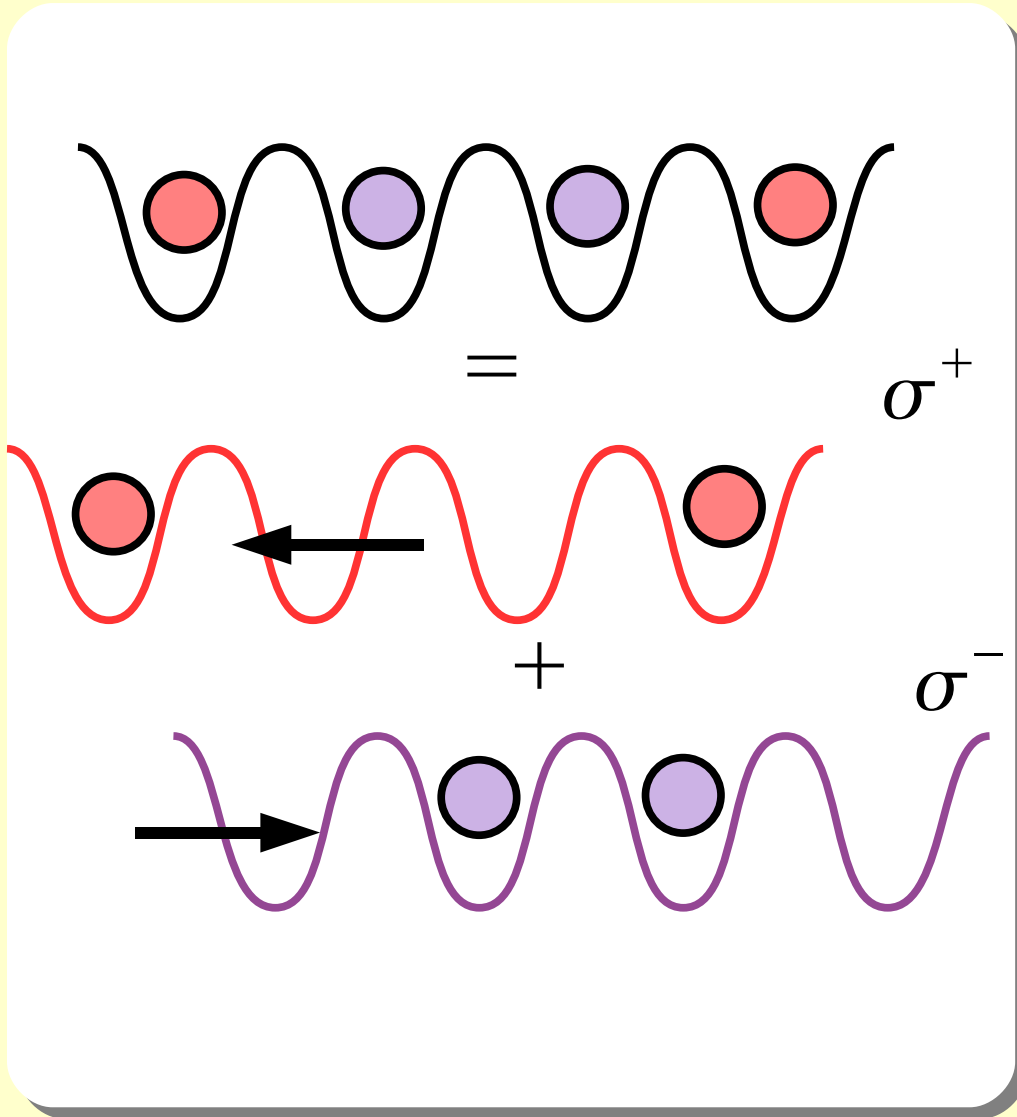
State-dependent lattices



- Not all atomic states react the same to different polarizations.
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- We can use this to create state-dependent lattices.
- We can apply a phase shift (birrefringence) to displace each lattice.
- The lattice has to be moved slowly, compared to the band gap.

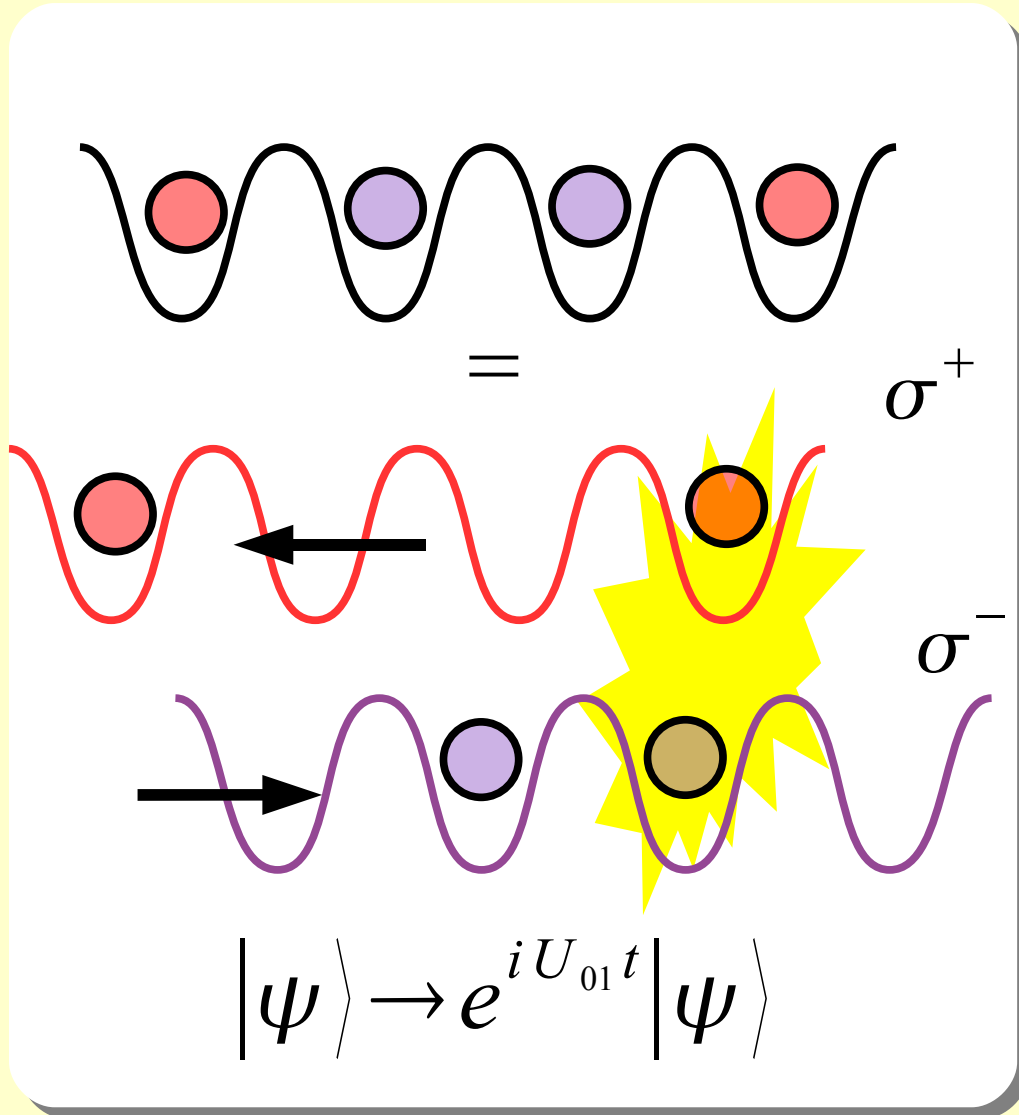
Quantum gates

Controlled collisions



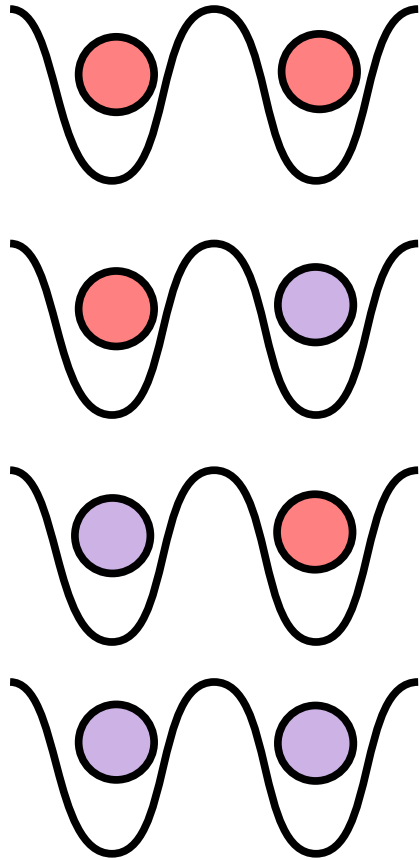
- We trap atoms in two states.
- Shift the lattices relative to each other.

Controlled collisions



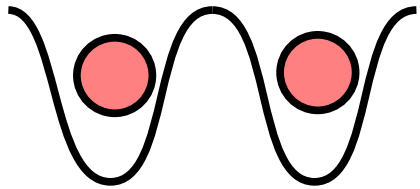
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- During the time collision takes place, they gain a phase.

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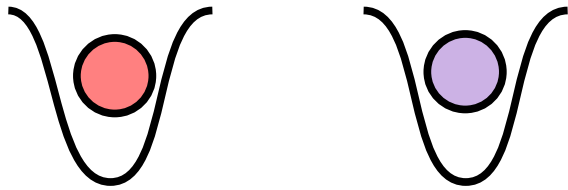


- We trap atoms in two states.
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- This phase implements a quantum gate.

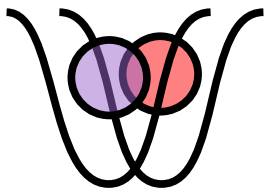
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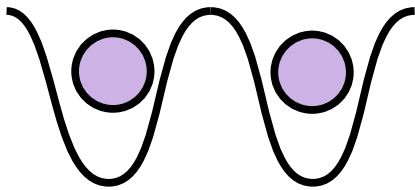
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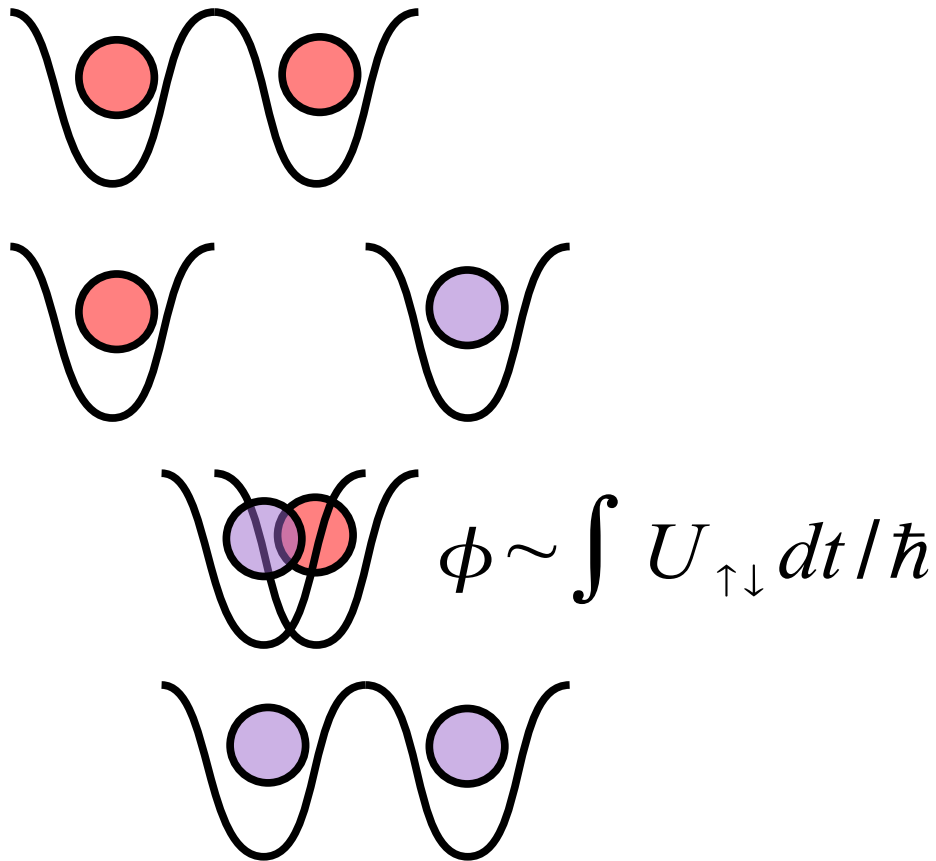
$$E \propto U_{\uparrow\downarrow}$$



$$E = 0$$

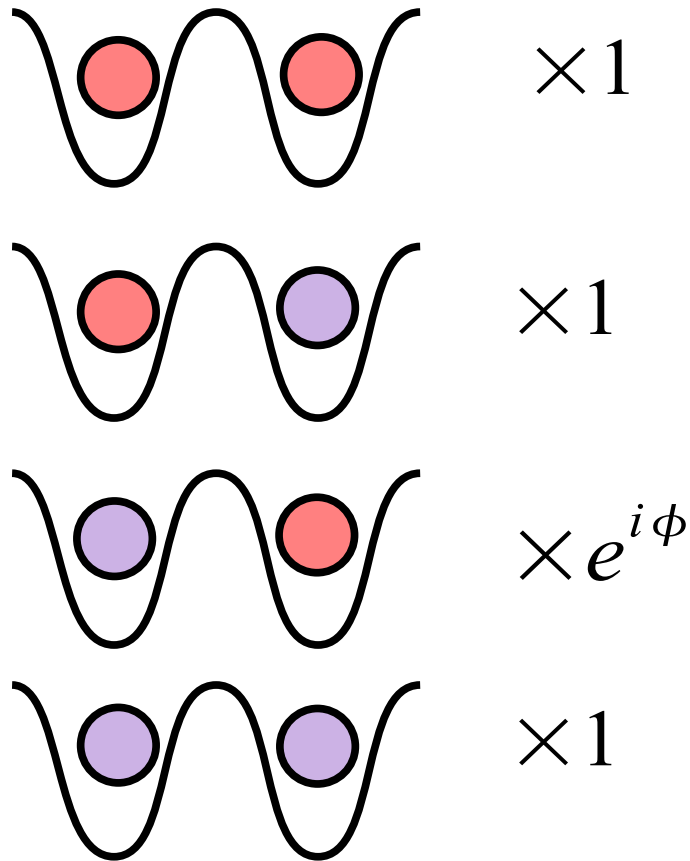
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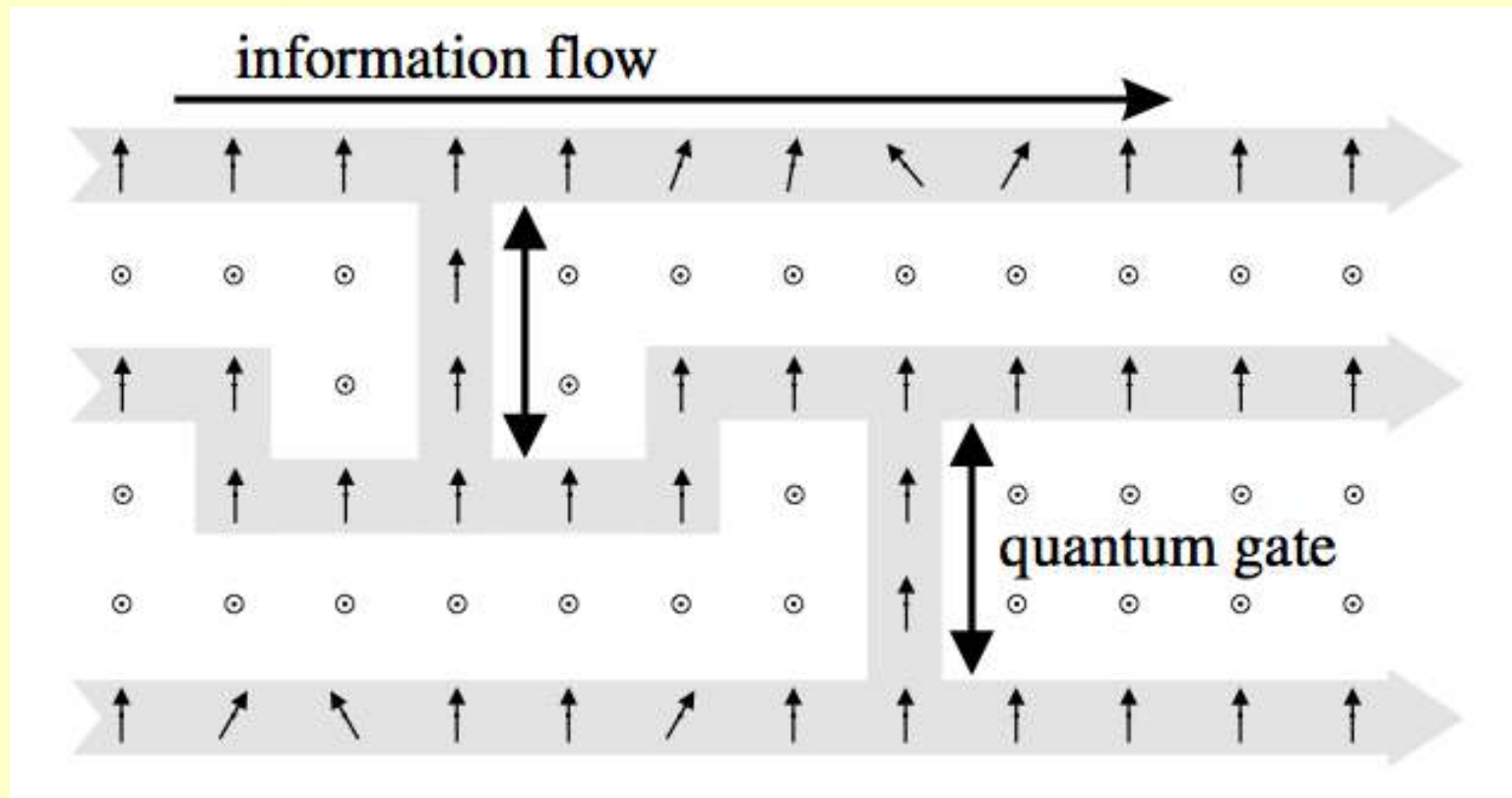


$$U_{gate} = e^{-iU_{01}t(\sigma_1^z - 1)(\sigma_2^z + 1)/4}$$

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Measurement based QC

Measurement based QC



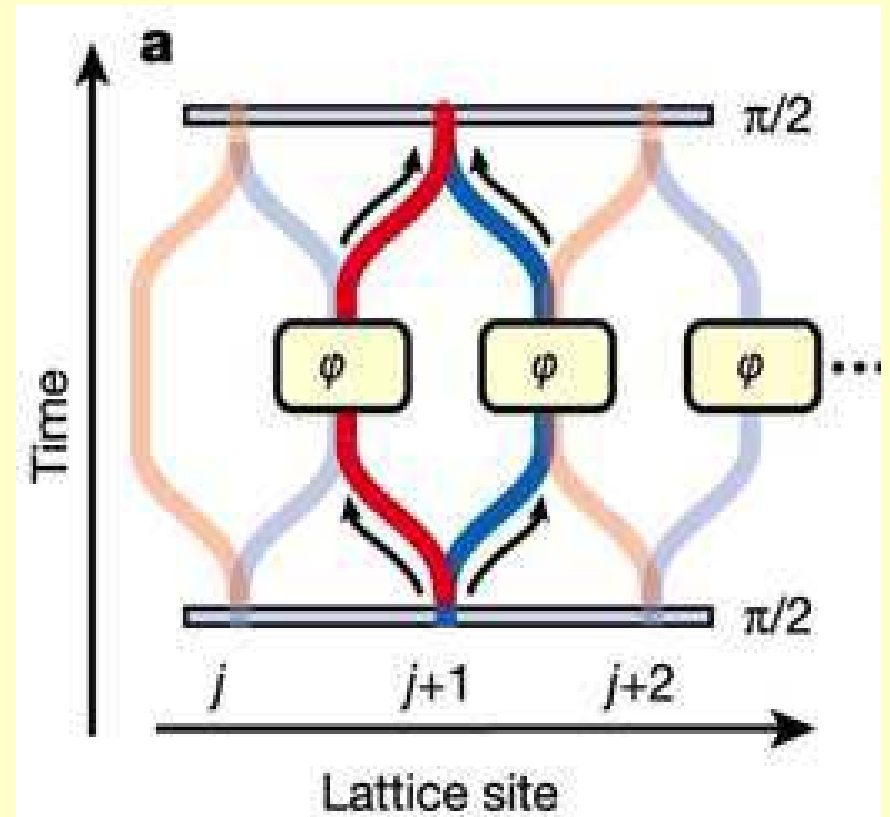
- It is possible to perform computations by measurements and local unitaries on a highly entangled state.
- This is called one-way quantum computing.

Cluster states

- In experiments this is done to **all atoms** at the same time.
- Massively parallelized quantum gate among nearest neighbors

$$U = \prod_{\langle i, j \rangle} e^{-i\phi[\sigma_i^z - 1][\sigma_j^z + 1]}$$

- Precisely what we need to make **cluster states**



$$U (|0\rangle + |1\rangle)^{\otimes N}$$

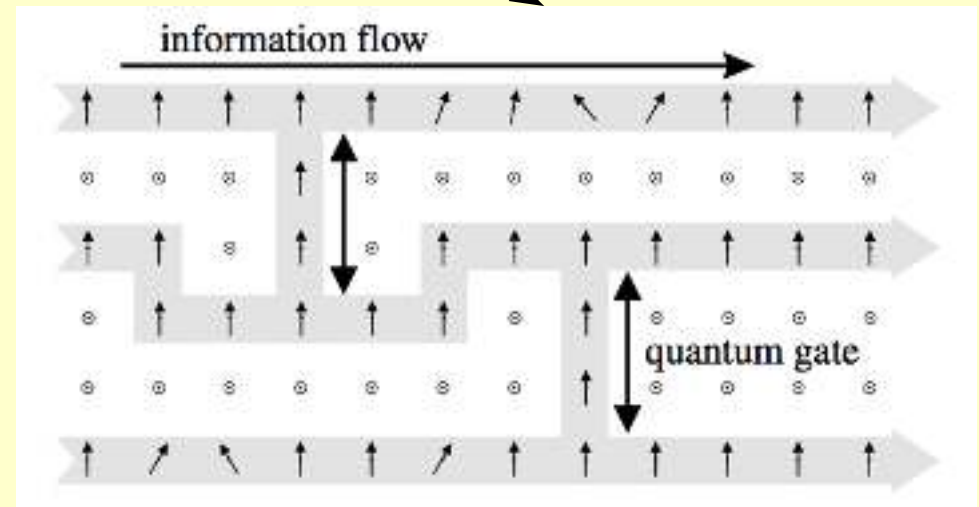
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Need individual addressing to make this useful



$$U (|0\rangle + |1\rangle)^{\otimes N}$$