

Other QC models

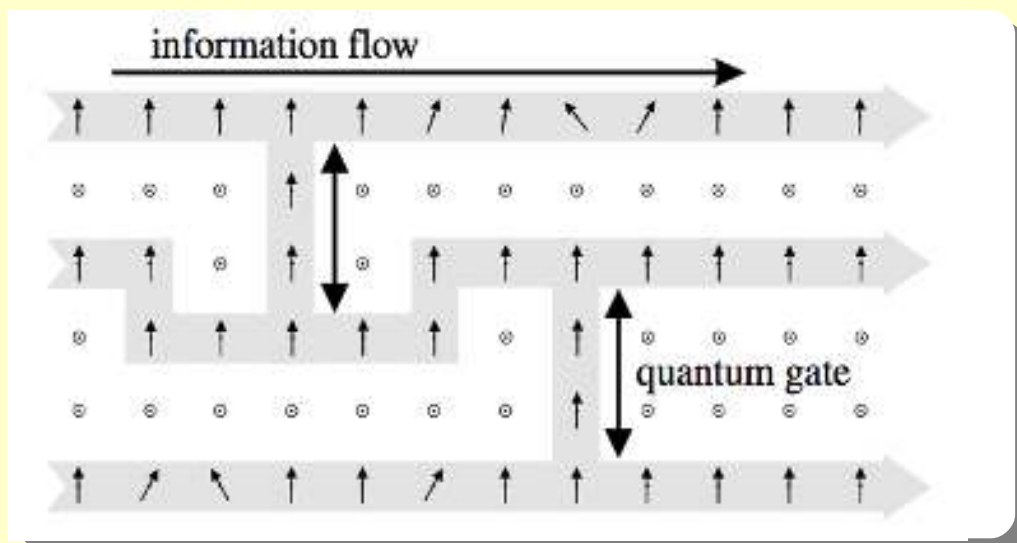
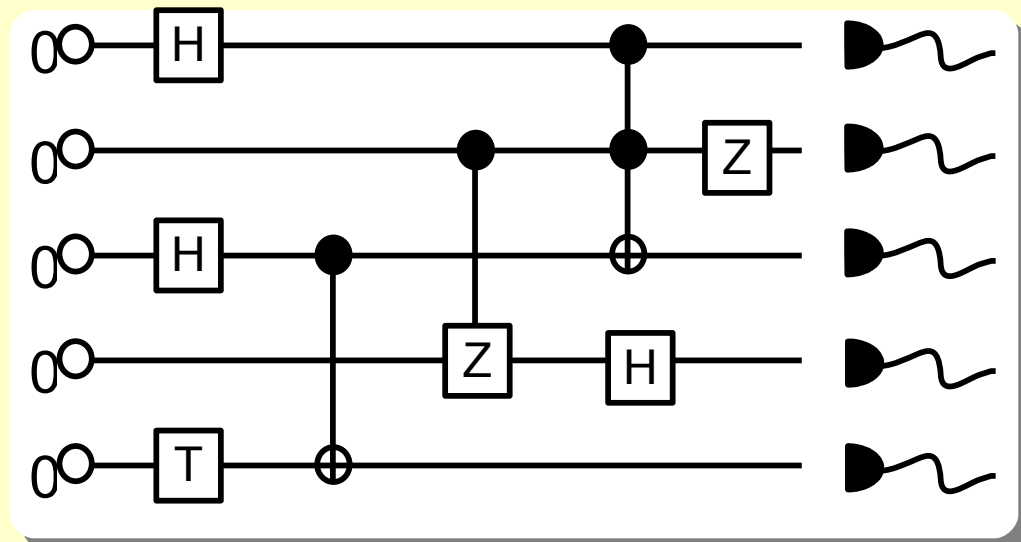
J. J. García-Ripoll
IFF, CSIC Madrid

(20-4-2009)

QC models so far

- A model based on a quantum register, combining local and universal two-qubit unitaries.
- Conceptually simple, parallelized operations.

Circuit based QC

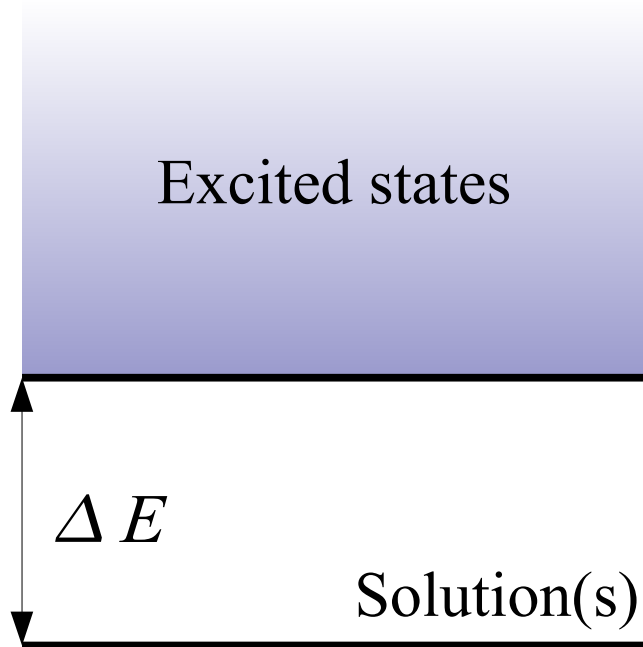


Measurement based QC

- Entangled state as a resource for QC.
- Only local operations and single-qubit measurements.
- Homogeneous, but same algorithms as circuit model

Adiabatic QC

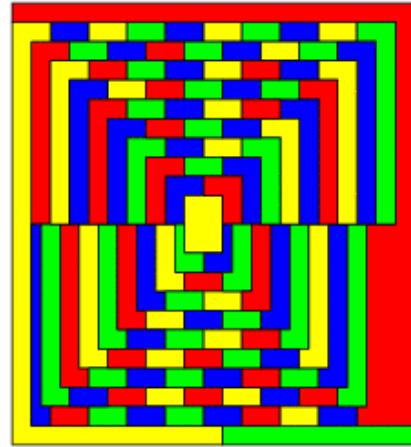
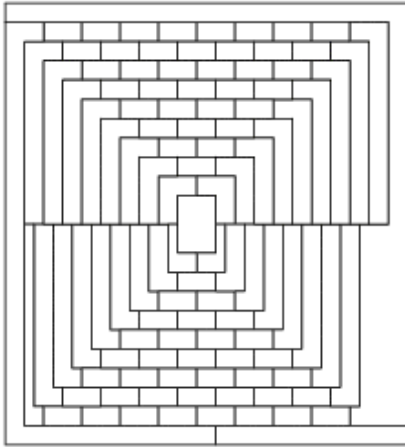
Adiabatic QC



$$(H_{prob} - E_{min})|\psi_{sol}\rangle = 0$$

- Express the outcome of a computation as a ground state.
- Design the appropriate interactions.
- Prepare adiabatically that ground state.
- Measure to find the outcome.

Utility?



Solve:

$$(-i_1) \vee i_3 \vee i_5 \wedge$$

$$i_2 \vee i_3 \vee (-i_5) \wedge$$

...

$$i_k \in \{0,1\}$$

Microchip optimization...

- Equivalent to quantum circuits.
- Many problems can be formulated as Hamilt.
 - Travelling salesman, nSAT, exact cover sudokus...
 - H_{prob} are simple boolean functions
- Problems *NP-complete*

How it works

$$H_0 = \sum_k \sigma_k^x$$

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



$$H(t) = t H_{prob} + (1 - t/T) H_0$$



$$H_{prob} = \sum_i f_i(\sigma_{i_1}^z, \sigma_{i_2}^z, \sigma_{i_3}^z)$$

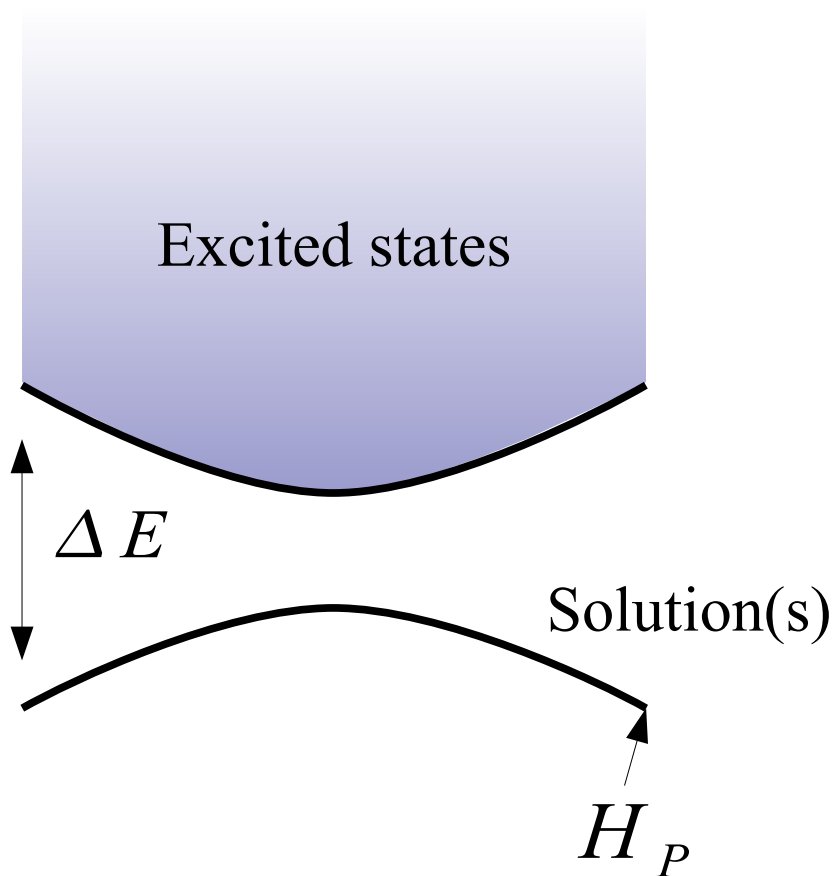
$$|\psi(T)\rangle \sim \sum_{\vec{s} \in Sol} |s_1, s_2 \dots s_N\rangle$$

- Begin with a state and H easy to prepare
- “Deform” adiabatically the effective interaction.
- The final H encodes the problem.
- We reach the final ground state if the instantaneous gap is large

$$\Delta E \gg \hbar/T$$

Example: Grover

Example: Grover



- We want to search a database, that is to solve
$$f(w)=1$$
- Our problem Hamiltonian can be simply

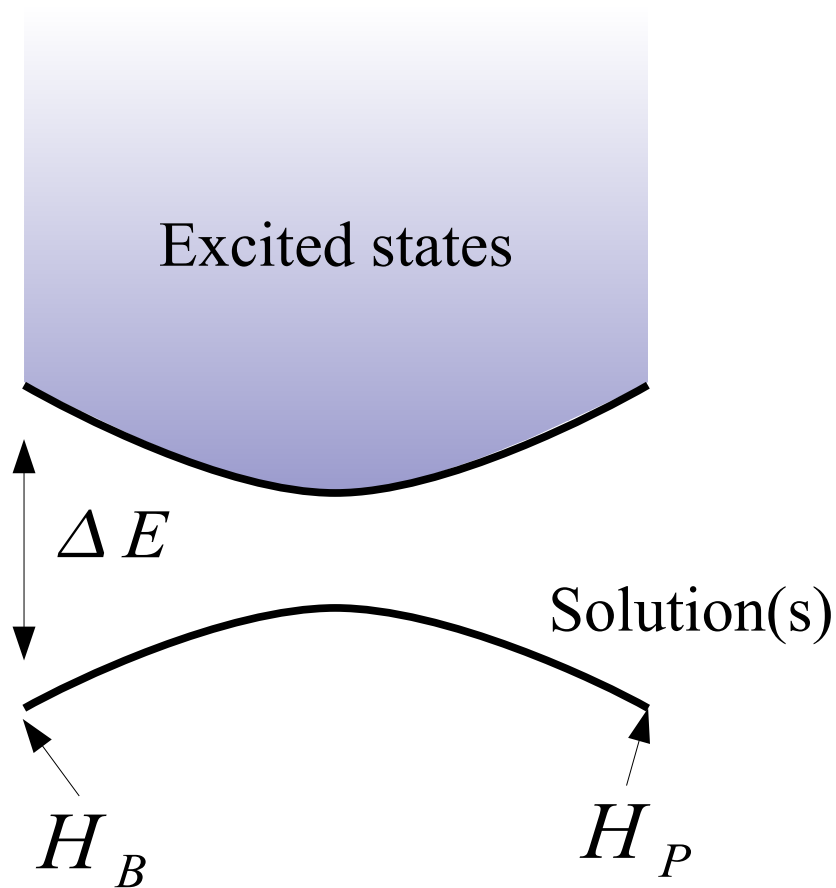
$$H_P = -|w\rangle\langle w|$$

with ground state $|w\rangle$

- Our Hamiltonian must end there

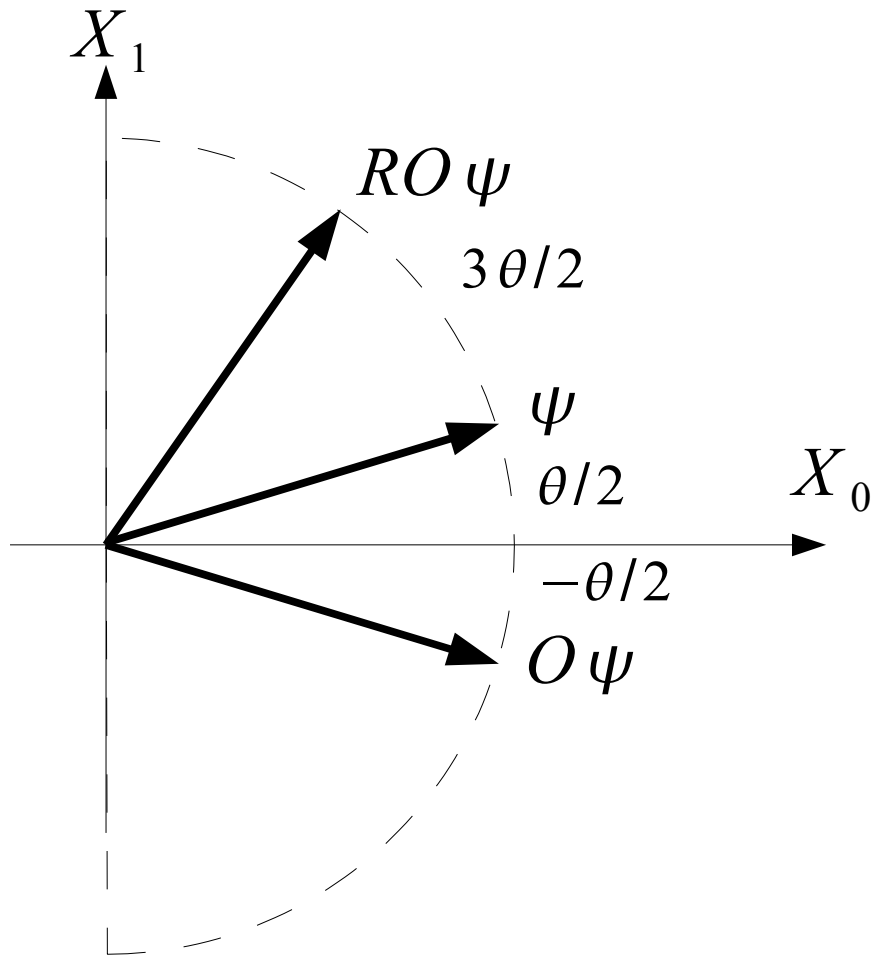
$$H(T) = H_P$$

Example: Grover



- The problem Hamiltonian has to be combined with a “mixing” term.
 - Allows transitions between different qubit states.
 - Easily prepared initial ground state.
- The interpolation among both may be as simple as
$$H(t) = \left(1 - \frac{t}{T}\right) H_B + \frac{t}{T} H_P$$
- This determines the gap and thus the alg. speed.

Mixing term



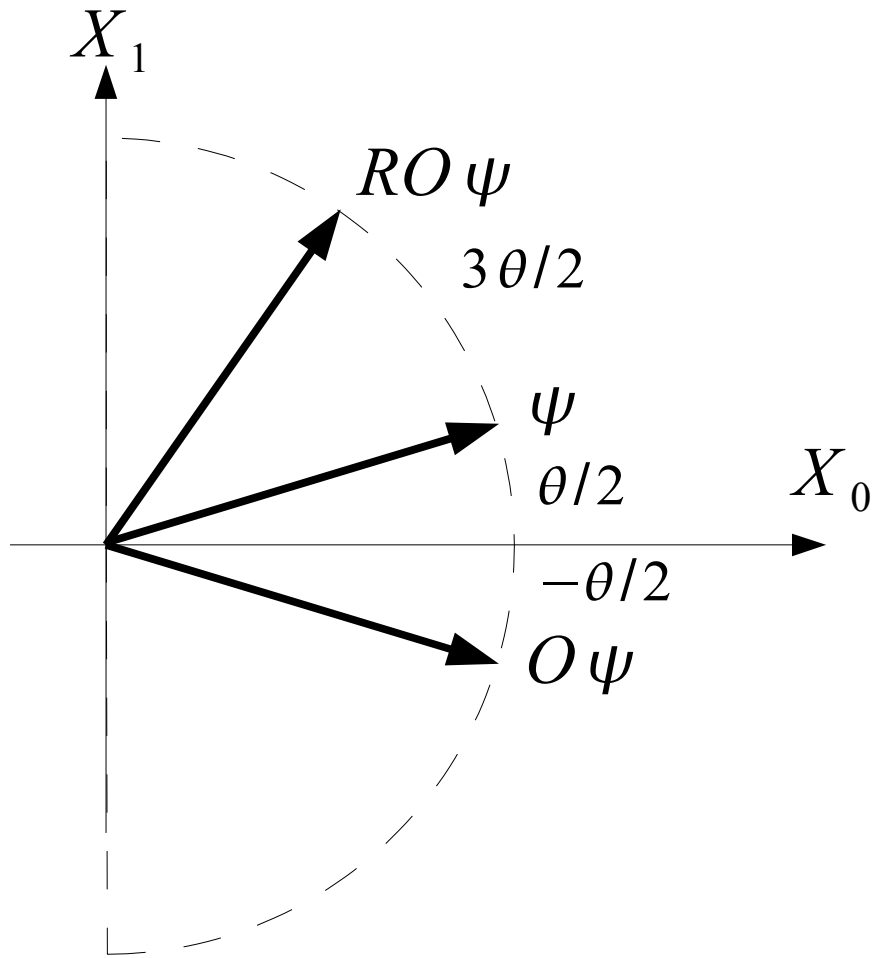
- We remind the fundamental operation in Grover

$$R \sim 2|\psi\rangle\langle\psi| - I$$

where

$$|\psi\rangle = \frac{1}{\sqrt{N}} \sum_{x=0}^N |x\rangle$$

Mixing term



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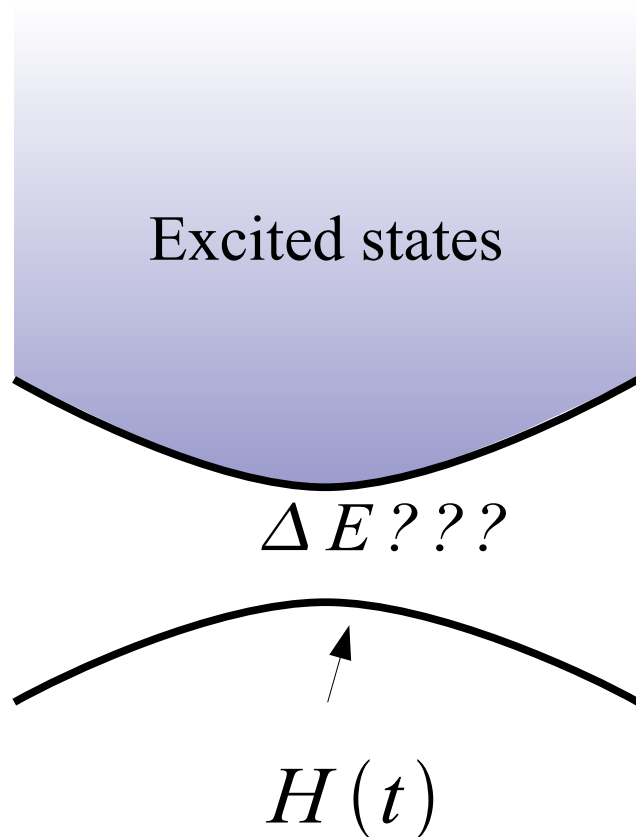
- We introduce thus a simple Hamiltonian

$$H_B = |\psi\rangle\langle\psi|$$

- Note that the initial state can be easily prepared.

$$|\psi\rangle \sim |+\rangle^{\otimes n}$$

Full Hamiltonian

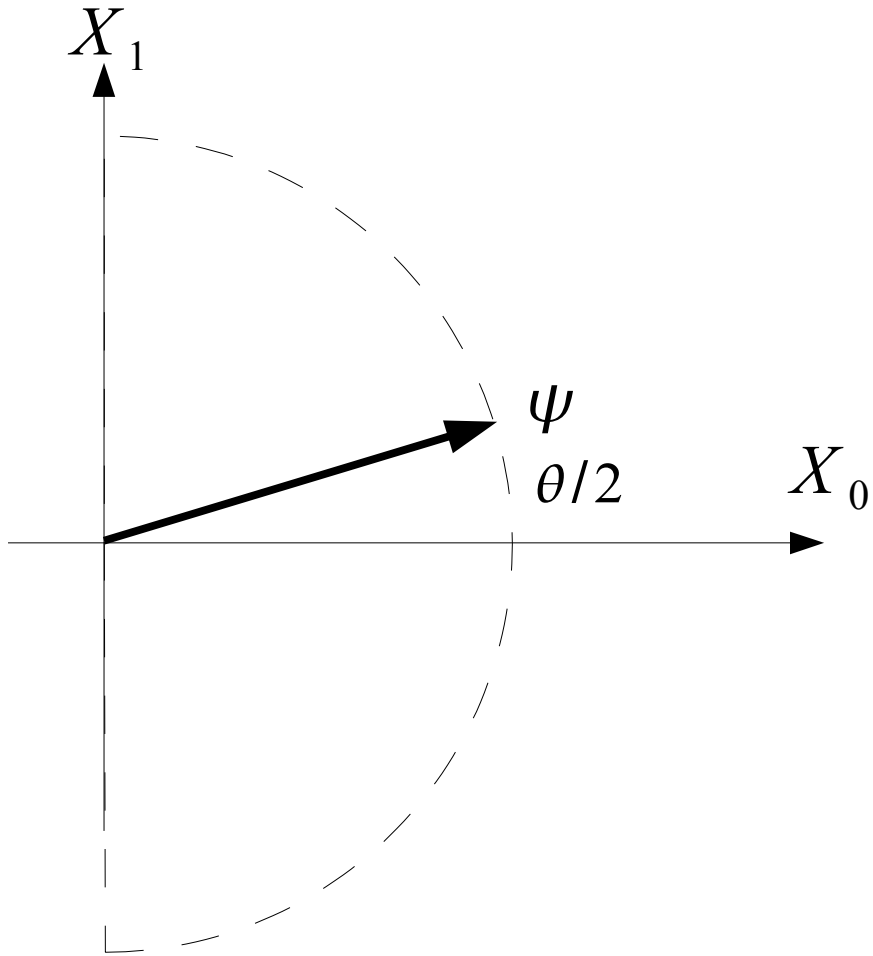


- We combine both terms

$$H = (1 - \lambda) \frac{1}{N} |\psi\rangle\langle\psi| - \lambda |w\rangle\langle w|$$

- Note that the energy is bound by $O(1)$
 - We could arbitrarily increase speed by enlarging $\|H\|$

Full Hamiltonian



- We reintroduce the idea from Grover

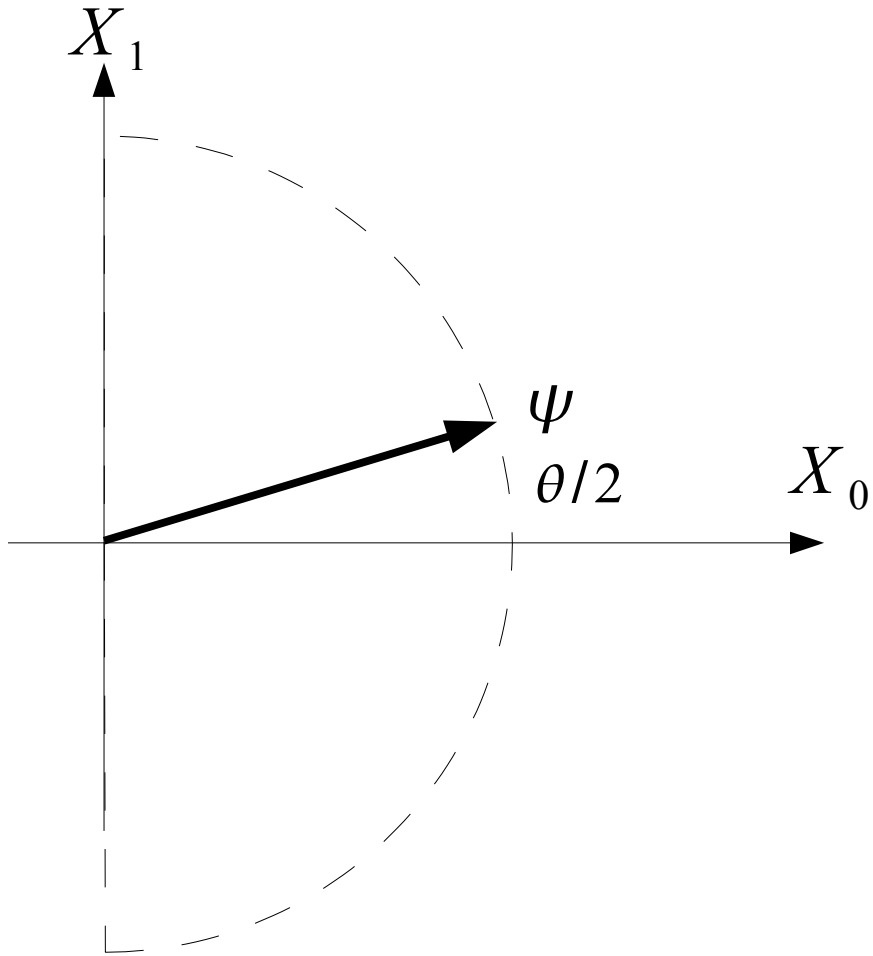
$$|\psi\rangle = \cos(\theta/2)|y\rangle + \sin(\theta/2)|w\rangle$$

so that

$$H_B = -\cos(\theta/2)^2|y\rangle\langle y| + \sin(\theta/2)^2|w\rangle\langle w| + \sin(\theta/2)\cos(\theta/2)[|w\rangle\langle y| + |y\rangle\langle w|]$$

$$H_P = -|w\rangle\langle w|$$

Full Hamiltonian



- Easier to read

$$H_B \sim -\frac{1}{2} - \cos(\theta) \sigma_z - \frac{1}{2} \sin(\theta) \sigma_x$$

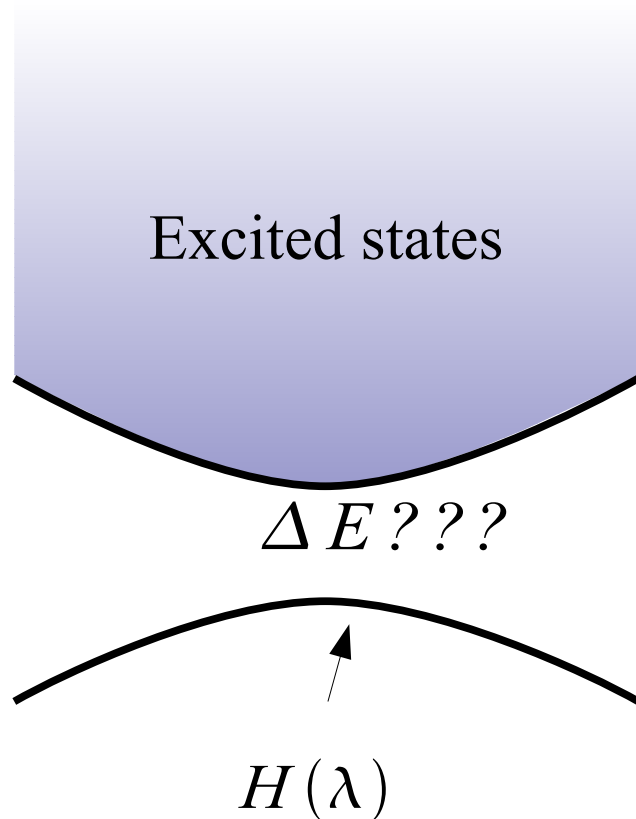
$$H_P \sim -\frac{1}{2} + \frac{1}{2} \sigma_z$$

using Pauli operators

$$\sigma_z = |y\rangle\langle y| - |w\rangle\langle w|$$

$$\sigma_x = |y\rangle\langle x| + |w\rangle\langle w|$$

Full Hamiltonian



- We can rewrite in the form

$$H = -\frac{1}{2} - \frac{1}{2} \vec{v} \cdot \vec{\sigma}$$

where

$$\vec{v} = [(1 - \lambda) \cos(\theta) - \lambda, (1 - \lambda) \sin(\theta)]$$

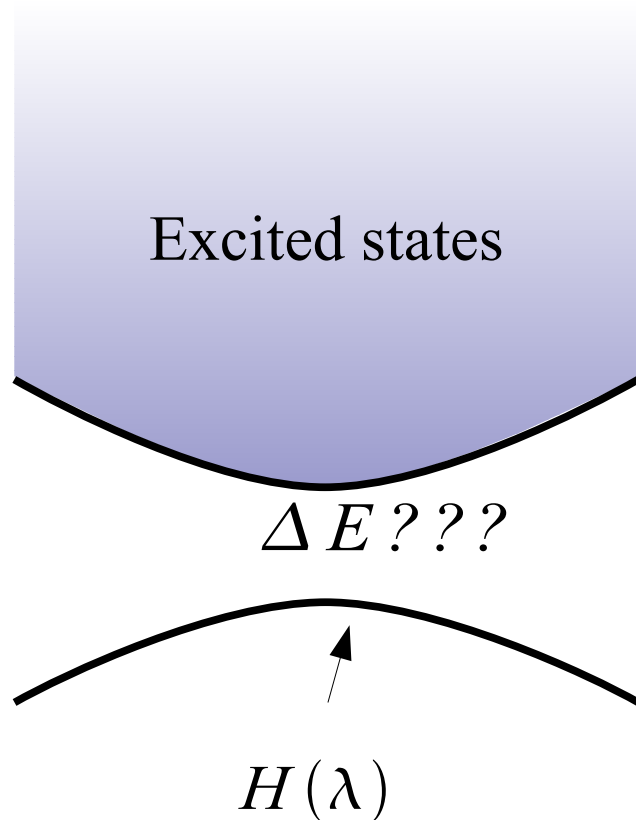
- The energy levels are

$$E = -\frac{1}{2} \pm \frac{1}{2} |\vec{v}|$$

with energy gap

$$\Delta E = |\vec{v}|$$

Full Hamiltonian



- From the expression

$$\vec{v} = [(1 - \lambda) \cos(\theta) - \lambda, (1 - \lambda) \sin(\theta)]$$

the gap is

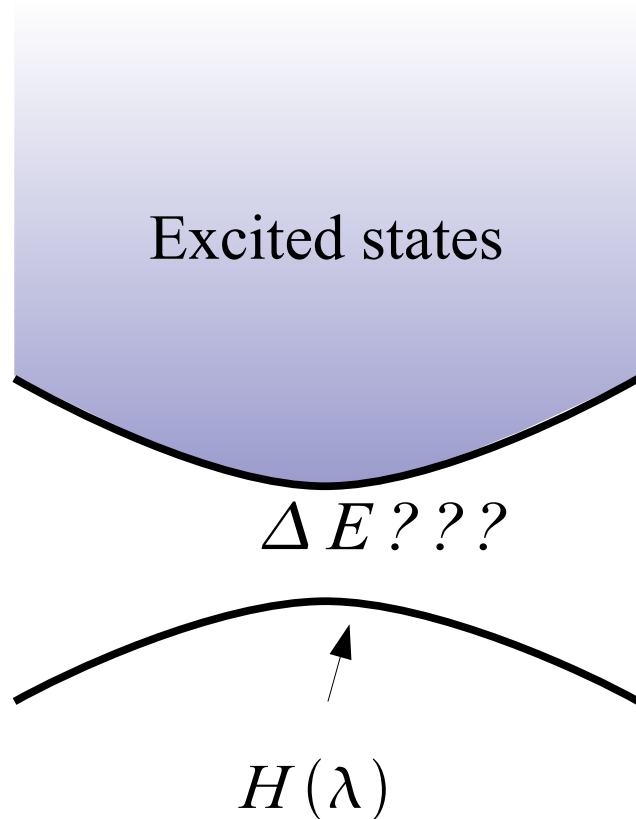
$$\Delta E = \sqrt{1 - 2\lambda(\lambda - 1)(\cos\theta + 1)}$$

- For our particular choice of state, the angle is

$$\cos(\theta/2)^2 = \frac{N-1}{N}$$

$$\cos(\theta) = 1 - \frac{2}{N}$$

Full Hamiltonian



- Summing up

$$\Delta E = \sqrt{1 - 2\lambda(\lambda - 1)(\cos\theta + 1)}$$

$$\cos(\theta) = 1 - \frac{2}{N}$$

$$\lambda_{min} = \frac{1}{2}$$

$$\Delta E_{min} = \frac{1}{\sqrt{N}}$$

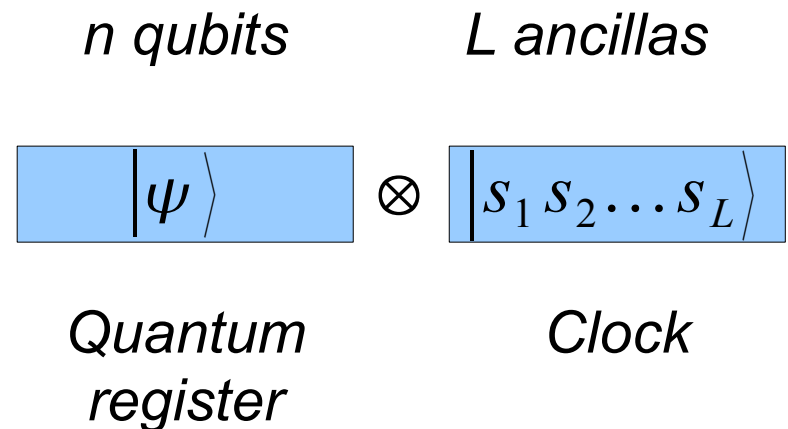
and time behaves like in Grover

$$T \geq O(\sqrt{N})$$

Why the interest?

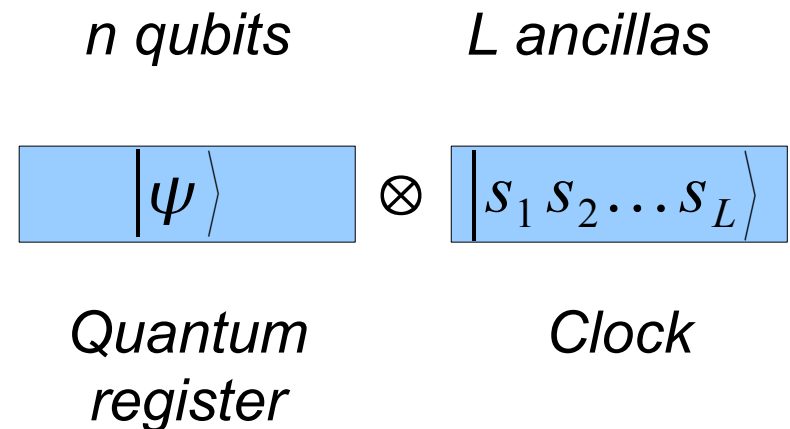
1) Universality

- Adiabatic quantum computation is as powerful as the circuit model.
- For every quantum circuit there exists an adiabatic Hamiltonian and viceversa.
- The overhead in the equivalence is polynomial
 - in time (adiab. resource)
 - in gates and ancillas (circuit resource)



1) Universality

- The proof in one sense uses Trotter decompositions
 - Approximate adiabatic evolution using quantum gates.
- On the other side we use
 - additional qubits to form a “clock”
 - a Hamiltonian that implements the unitaries
 - a Hamiltonian that forces the clock states to be consistent

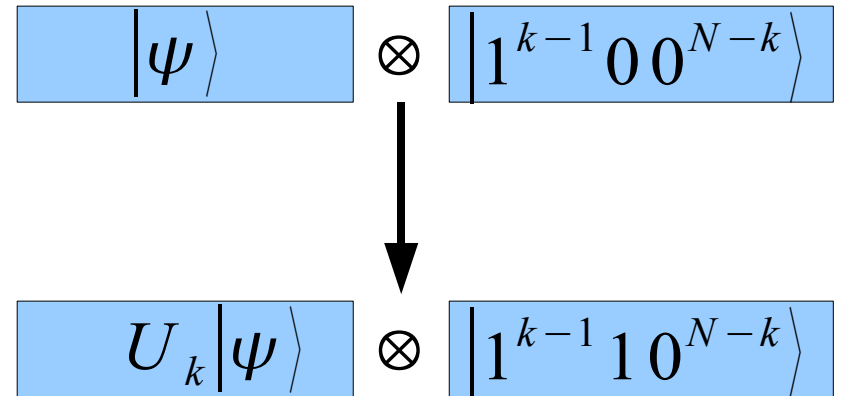


1) Universality

- The unitary Hamiltonian imposes that transitions only happen for the right clock states

$$\begin{aligned} H_k = & I \otimes |100\rangle\langle 100|_{k-1,k,k+1} \\ & + I \otimes |110\rangle\langle 110|_{k-1,k,k+1} \\ & + U_k \otimes |110\rangle\langle 100|_{k-1,k,k+1} \\ & + U_k^+ \otimes |100\rangle\langle 110|_{k-1,k,k+1} \end{aligned}$$

k-th circuit step



1) Universality

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- Another term forbids wrong clock states

$$H_{clock} = \sum_k |01\rangle\langle 01|_{kk+1}$$

consistent clock

$$\boxed{|\psi\rangle} \otimes \boxed{|1\dots 01\dots\rangle}$$

NO!

1) Universality

- Another term imposes the initial state at zero clock

$$H_{init} = \left[\sum_i |1\rangle\langle 1|_i \right] \otimes |0\rangle\langle 0|_0^{clock}$$

- And we have to impose initially the clock is zero

$$H_{clock0} = \sum_k |1\rangle\langle 1|_k^{clock}$$

initial state



The diagram illustrates the initial state as a tensor product of two states. It consists of two light blue rectangular boxes, each containing the quantum state $|0 \dots 0\rangle$. The two boxes are separated by a tensor product symbol \otimes .

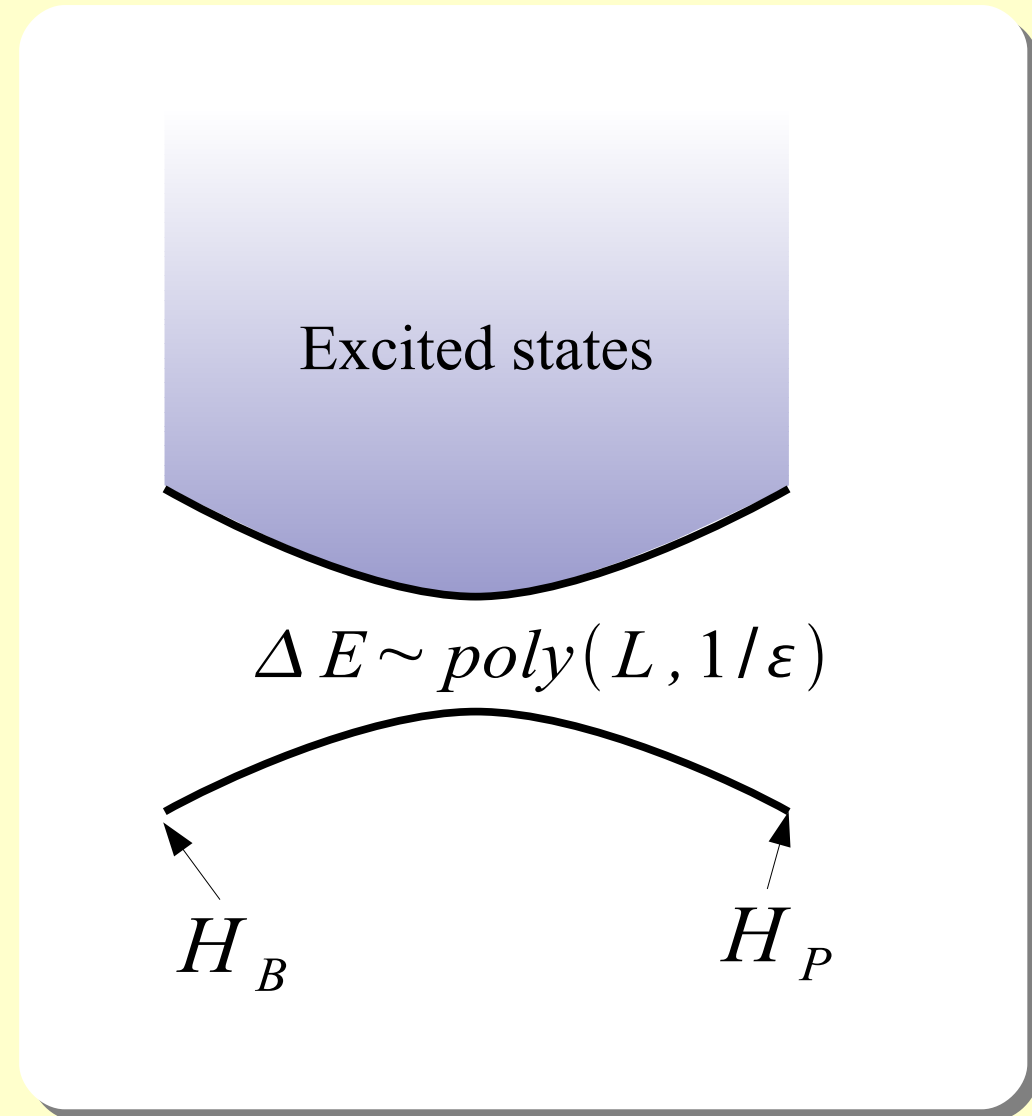
1) Universality

- Put together

$$H_B = H_{clock0} + H_{init} + H_{clock}$$

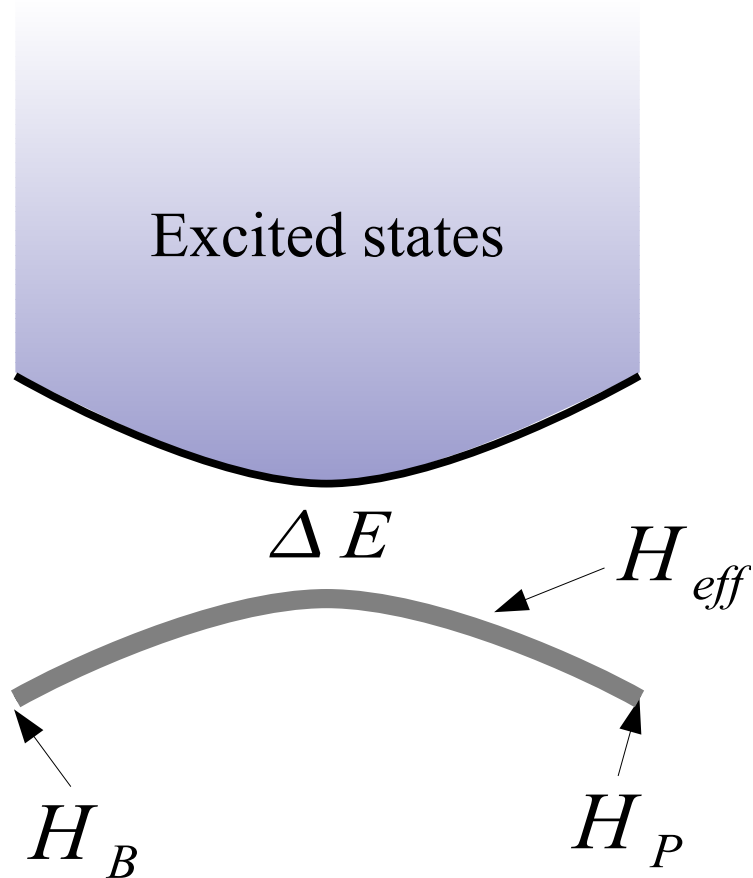
$$H_P = \sum_k H_k + H_{init} + H_{clock}$$

- The merit of the proof is to show that the gap is polynomial in the problem size, L , and inverse tolerance
 - Time increases with the number of gates still!



2) Robustness

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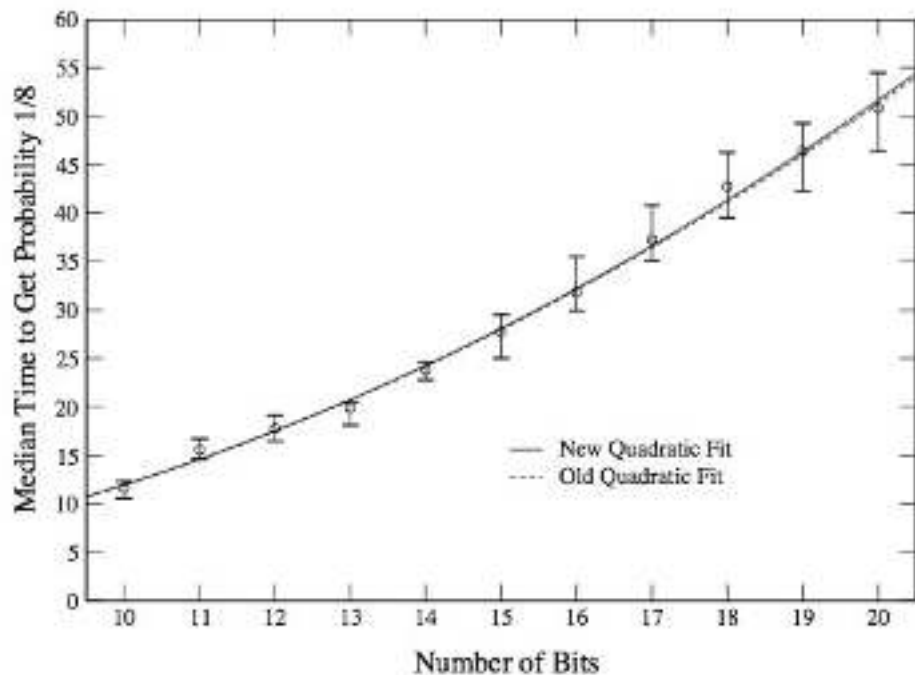
- Remember perturbation theory:
 - We have two separate regions in spectrum.
 - If perturbations are small, they remain disconnected.
 - At most a new effective Hamiltonian for the ground state manifold
$$H_{eff} = (1 - \lambda) H_B + \lambda H_P + \epsilon H_{eff}$$
 - At most mixing of ground states.

Real life

Local Hamiltonians

- What mixing Hamiltonians are experimentally realizable?

Local Hamiltonians



- What mixing Hamiltonians are experimentally realizable?
- We have use a very nonlocal one

$$H_B = |\psi\rangle\langle\psi| = \frac{1}{N} \sum_{x, y=0}^N |x\rangle\langle y|$$

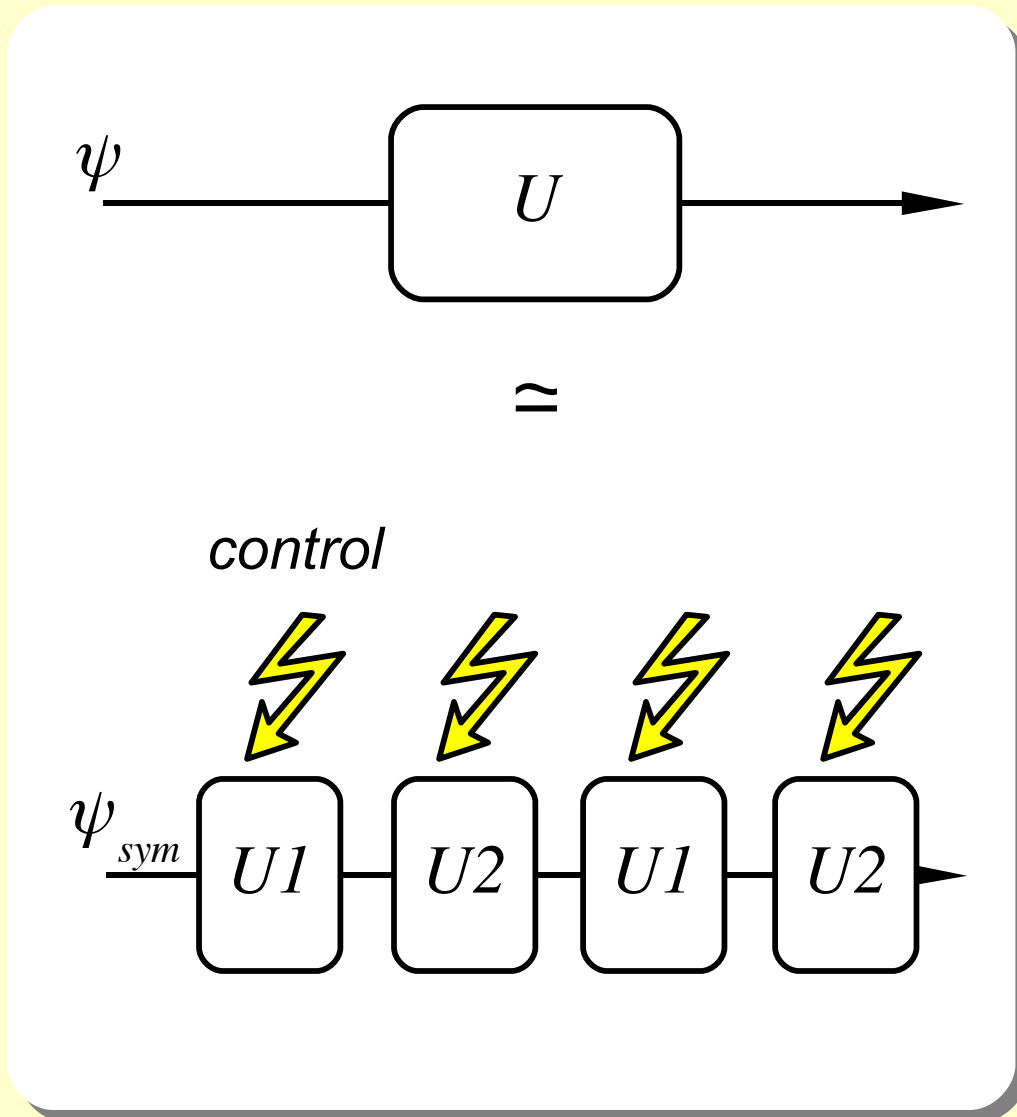
- An alternative would be

$$H_B = \sum_{i=0}^n \sigma_i^x$$

- Computing gap is difficult
- Little numerical evidence

Practical implementations

Simulated implementation

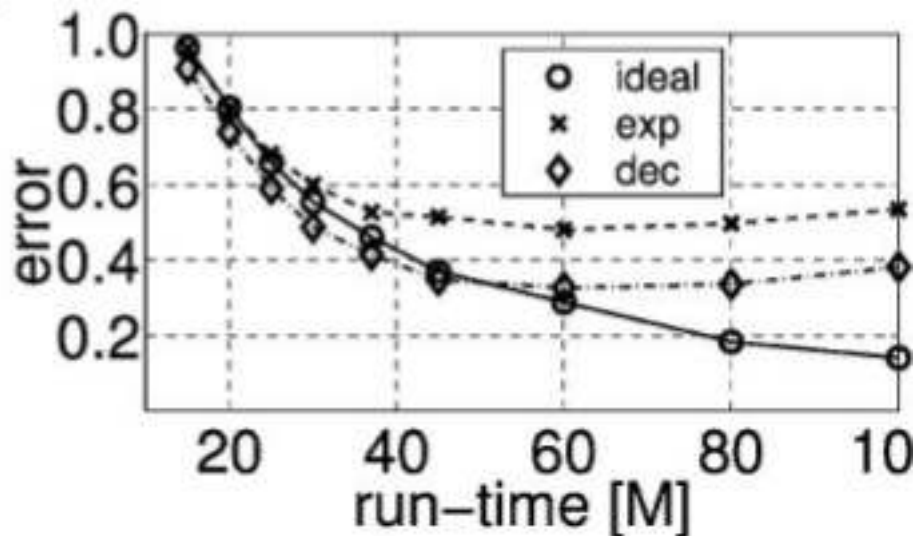


- An adiabatic evolution can be simulated using two alternating Hamiltonians

$$U(T) \sim \prod_n \exp \left\{ i \left[(1 - \lambda_n) H_B + \lambda_n H_P \right] \right\} \\ \sim \prod_n \exp \left\{ i (1 - \lambda_n) H_B \right\} \times \\ \times \exp \left\{ i \lambda_n H_P \right\}$$

- This introduces additional errors, but we expect the effective “gap” will suppress them.

NMR adiabatic algorithm



- Molecule: H F C Br2
- Effective spin-spin interaction between H, F and C nuclei

$$H = \sum_{i,j} J_{i,j} \sigma_i^z \sigma_j^z + \sum_i \left[B \sigma_i^z + \vec{B}(t) \cdot \vec{\sigma} \right]$$

- With a suitable modulation of the transverse mag. field, we can simulate both the problem

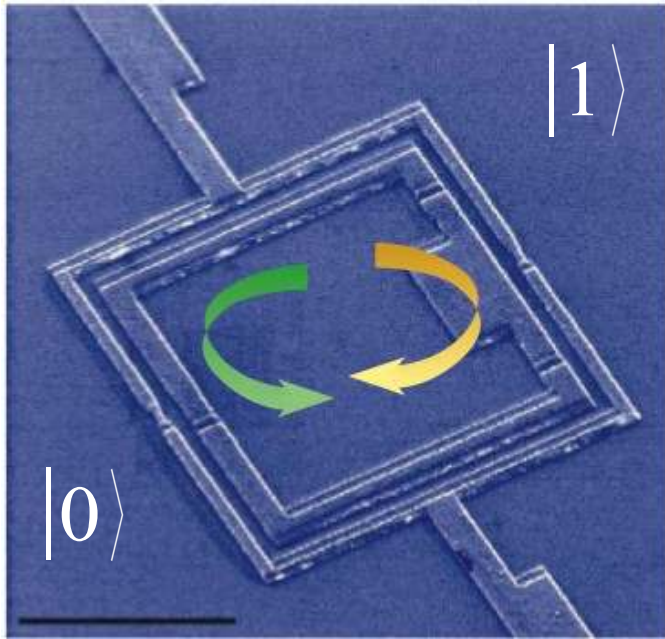
$$H_P = \sum_{i,j} f(\sigma_i^z, \sigma_j^z)$$

and mixing Hamiltonians

$$H_B = \sum_i \sigma_i^x$$

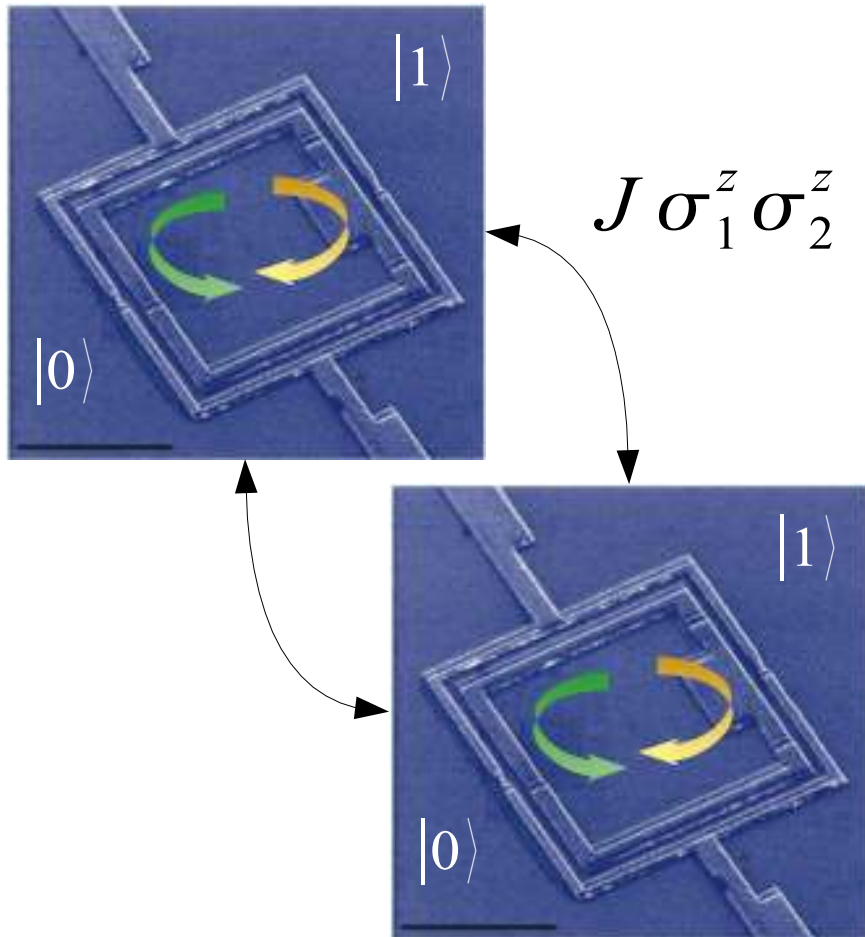
“Real” adiabatic computation

Superconducting circuits



- SQUIDS are superconducting loops.
- They can host long lived currents.
- The sense of a current can implement a qubit state

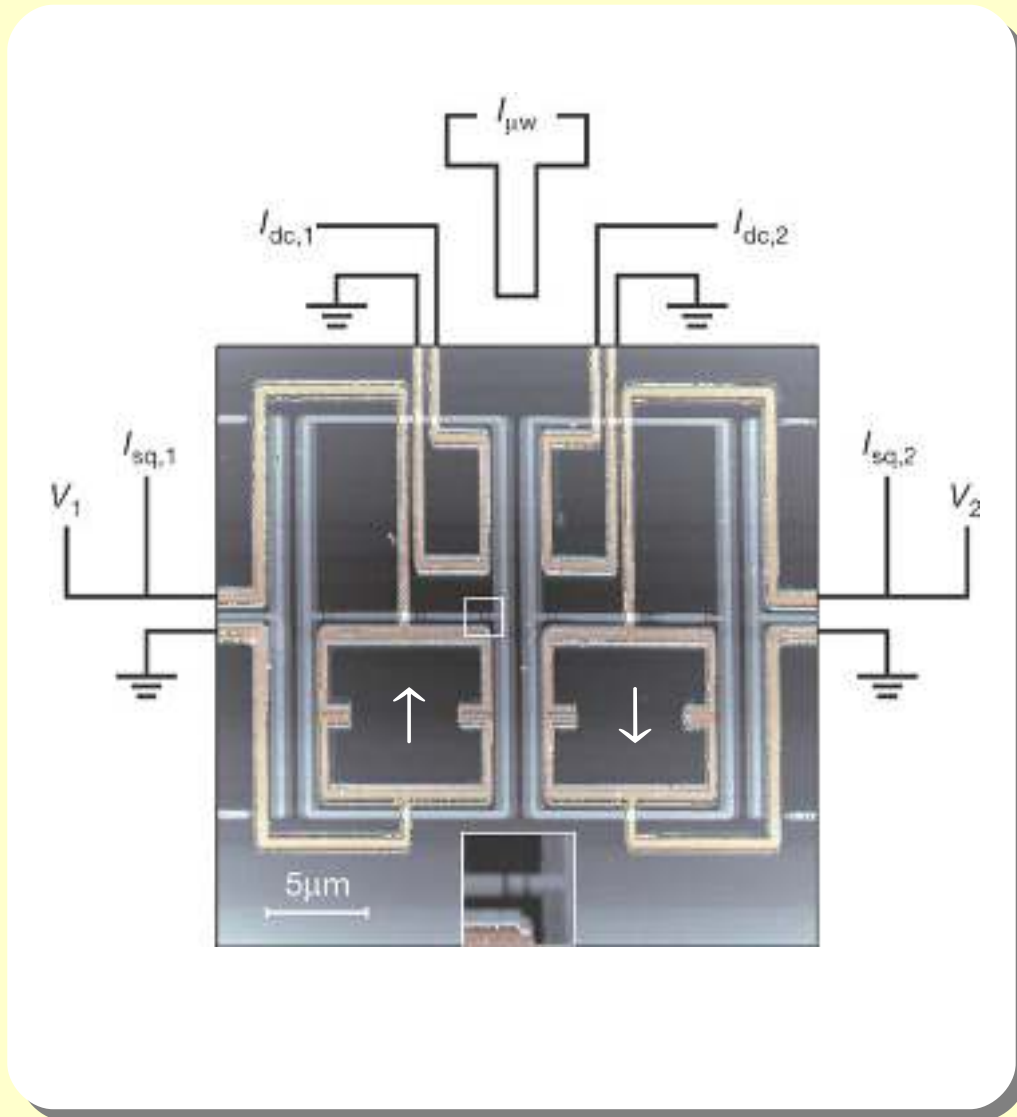
Superconducting circuits



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- These currents behave like small **magnets** giving rise to a magnetic interaction

$$H \sim \sum_{i,j} J_{i,j} \sigma_1^z \sigma_2^z + \sum h_i^{ext} \sigma_i^x$$

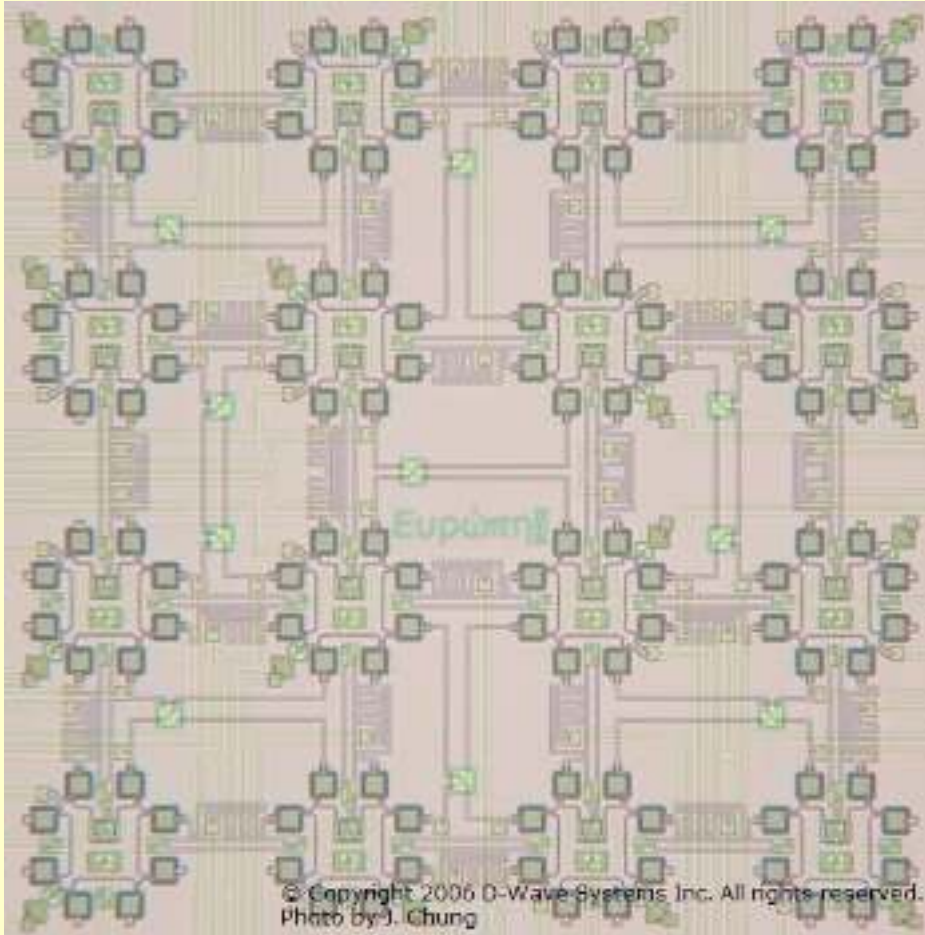
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Enter D-Wave



- 4x4 superconducting qubits of flux or phase type.
- Claims:
 - Adjustable interactions.
 - Implements adiabatic quantum algorithms
 - Should solve optimization problems.

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SCIENTIFIC AMERICAN

[News](#) - February 13, 2007

First "Commercial" Quantum Computer Solves Sudoku Puzzles

Quantum computing company banks on a long-shot form of quantum computing

The New York Times
nytimes.com

April 8, 2007

SLIPSTREAM

A Giant Leap Forward in Computing?

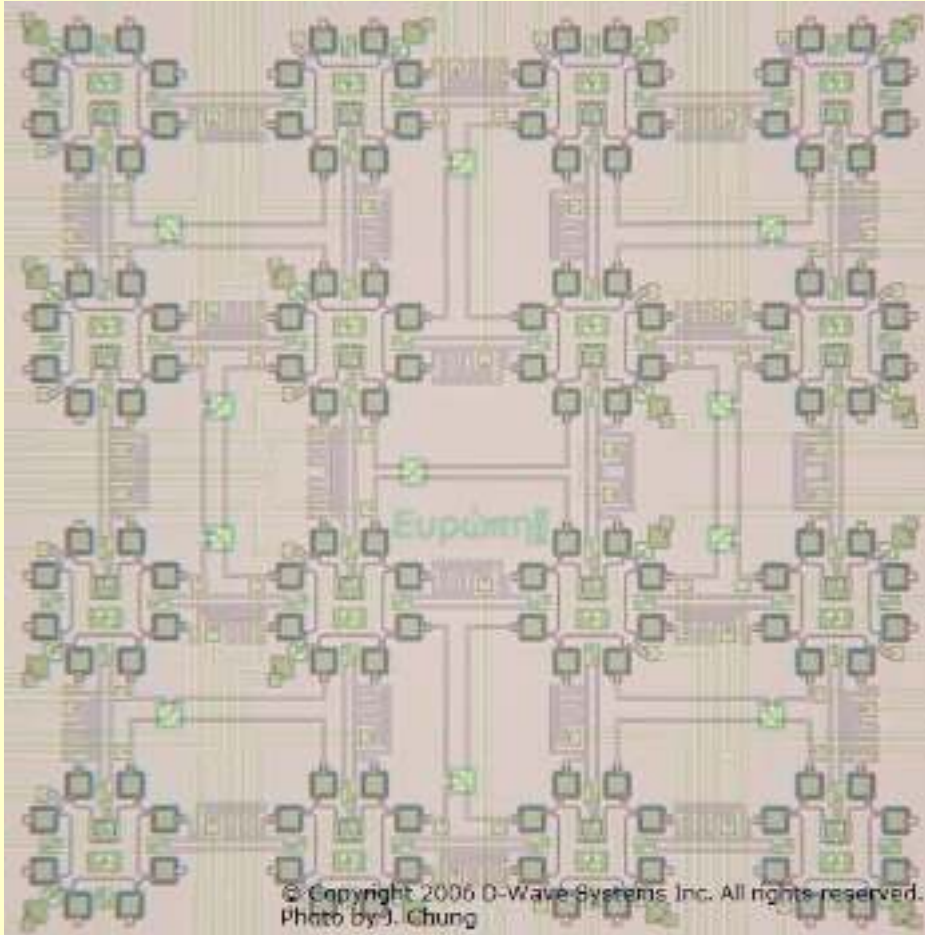
 **NEWS.COM**

<http://www.news.com/>

Start-up demos quantum computer

By Michael Kanellos

Problems



- Adiabatic algorithm requires coherence
 - Otherwise annealing
- In D-Wave qubits are closer to classical bits with noise.
- But we cannot discard it as a possibility!