

Title: More on adiabatic quantum computation - Part 2

Date: Feb 11, 2006 09:50 AM

URL: <http://pirsa.org/06020023>

Abstract:

$$\text{Grover } + (z) = \begin{cases} -1 & z = w \\ 0 & z \neq w \end{cases}$$

$$0 \leq z \leq N-1$$

Find w

Classically N calls

Q.M. $N^{\frac{1}{2}}$ calls \leftarrow Best possible

$$H_p = H_w = E(|-\rangle\langle w|)$$

Can apply H_p but do not know w

$$H(t) = H_0(t) + H_w$$

e.g. You design this

$$\text{Want } |\psi(t)\rangle = |w\rangle$$

$$T \geq \frac{\sqrt{N}}{2E}$$

E.F.
S. Gutmann

4.

● Proof Use $H_w = E|\omega\rangle\langle\omega|$

$$H = H_0(t) + H_w$$

$$t=0 \quad |\text{initial}\rangle$$

$$t=T \quad |\omega\rangle \quad \text{want}$$

$$i\frac{d}{dt}|\psi_w, t\rangle = [H_0(t) + H_w]|\psi_w, t\rangle$$

$$|\psi_{w,0}\rangle = |\text{initial}\rangle; |\psi_{w,T}\rangle = |\omega\rangle$$

● Reference State $|\psi, t\rangle$ evolves with just $H_0(t)$

$$i\frac{d}{dt}|\psi, t\rangle = H_0(t)|\psi, t\rangle$$

$$|\psi, 0\rangle = |\text{init.}\rangle$$

$$\sum_w ||\psi_{w,T}\rangle - |\psi, T\rangle|^2 = \sum_w |\langle\omega| - \langle\psi, T|\omega\rangle|^2$$

$$= 2N - \sum_w [\langle\omega|\psi, T\rangle + \langle\psi, T|\omega\rangle]$$

$$\geq N$$

$$i \frac{d}{dt} |\psi(t)\rangle = H(t) |\psi(t)\rangle$$

$h(\cdot)$

$$i \frac{d}{dt} |\psi(t)\rangle = H(t) |\psi(t)\rangle$$
$$|\psi(T)\rangle$$

$\epsilon_{\text{h.r.}}$

$h(\epsilon)$

4.

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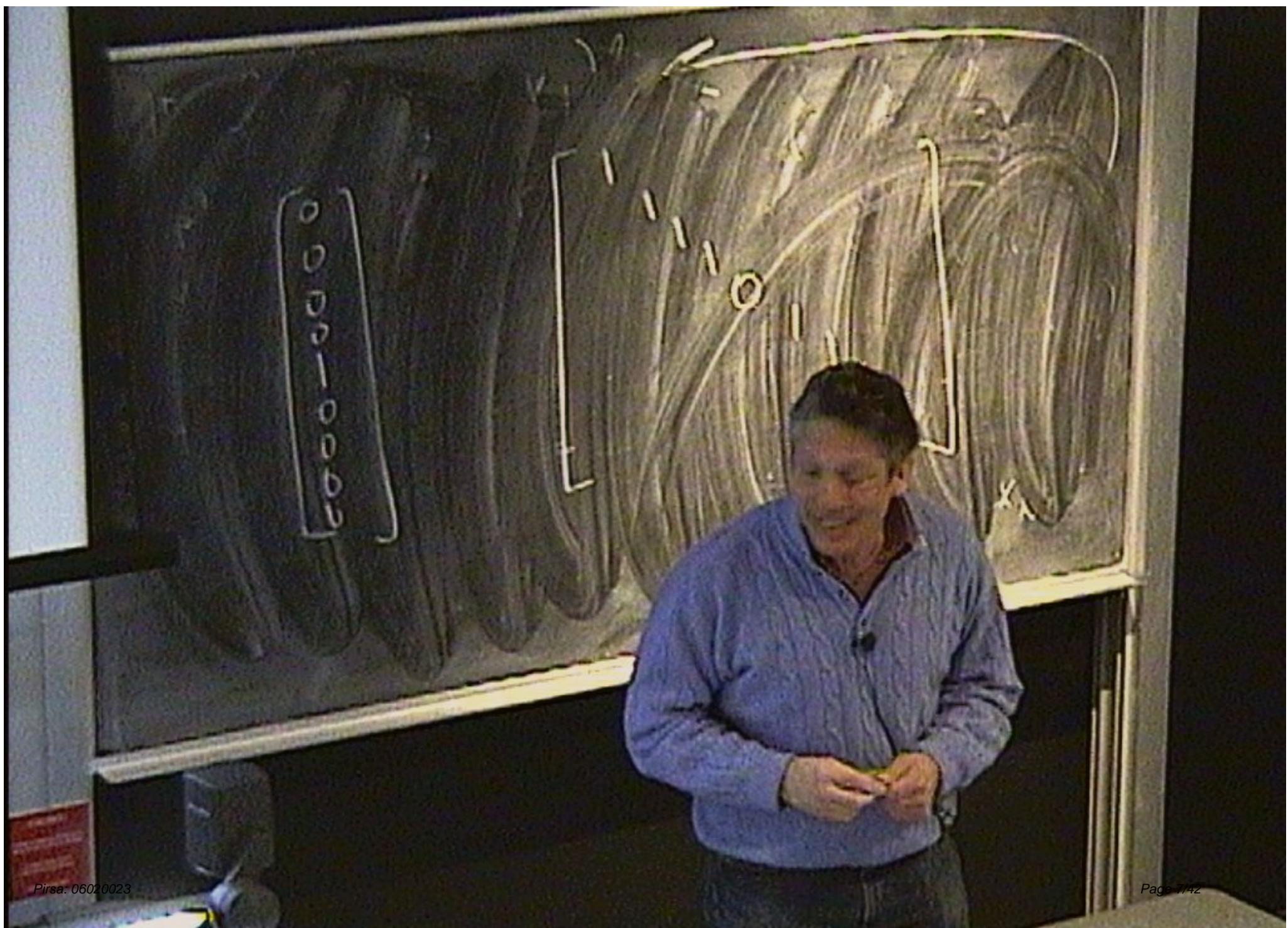
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$$i \frac{d}{dt} |\psi_{\omega}, t\rangle = [H_0(t) + H_w] |\psi_{\omega}, t\rangle$$

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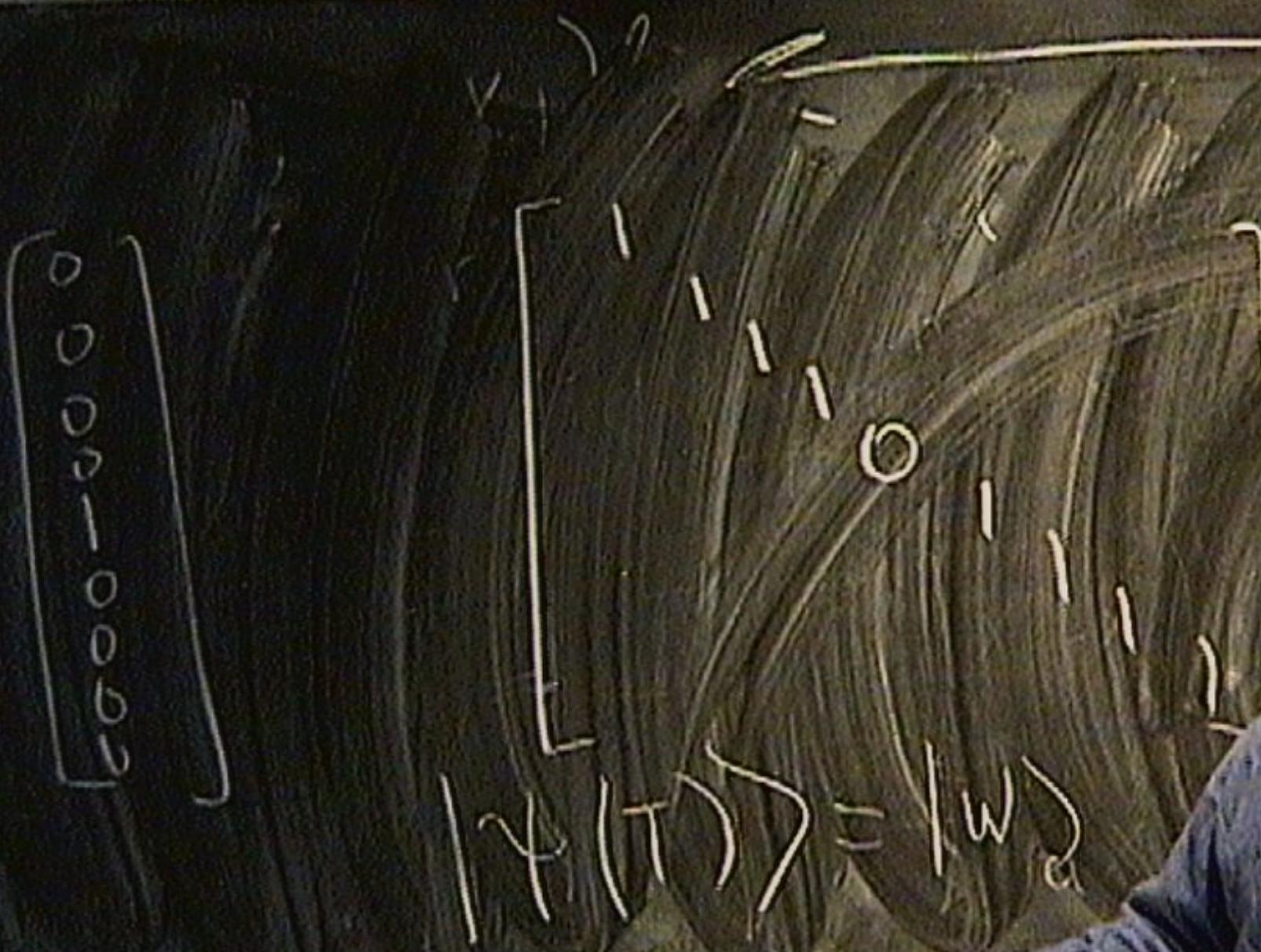
$$\sum_{\omega} | |\psi_{\omega, T}\rangle - |\psi, T\rangle |^2 = \sum_{\omega} | |\omega\rangle - |\psi, T\rangle |^2$$

$$= 2N - \sum_{\omega} [\langle \omega | \psi, T \rangle + \langle \psi, T | \omega \rangle]$$

$$\geq N$$

$$\frac{i \frac{d}{dt} |\psi(t)\rangle = H(t) |\psi(t)\rangle}{|\psi(t)\rangle =}$$

$h(t)$



● Proof Use $H_w = E|\omega\rangle\langle\omega|$

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$$= 2N - \sum_{\omega} [\omega|\psi, T\rangle + \langle\psi, T|\omega\rangle]$$

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$$\geq N$$

5.

$$\bullet \frac{d}{dt} |\langle \psi_\omega, t \rangle - \langle \psi_i, t \rangle|^2 = -2 \operatorname{Re} \frac{d}{dt} \langle \psi_\omega, t | \psi_i, t \rangle$$

$$= 2 \operatorname{Im} \langle \psi_\omega, t | H_\omega | \psi_i, t \rangle$$

$$\leq 2 \|H_\omega\| |\psi_i, t\rangle$$

Now \sum_{ω} : $\sum_{\omega} \in \langle \omega \rangle \langle \omega \rangle = E$

$$\frac{d}{dt} \sum_{\omega} |\langle \psi_\omega, t \rangle - \langle \psi_i, t \rangle|^2 \leq 2E N^{1/2}$$

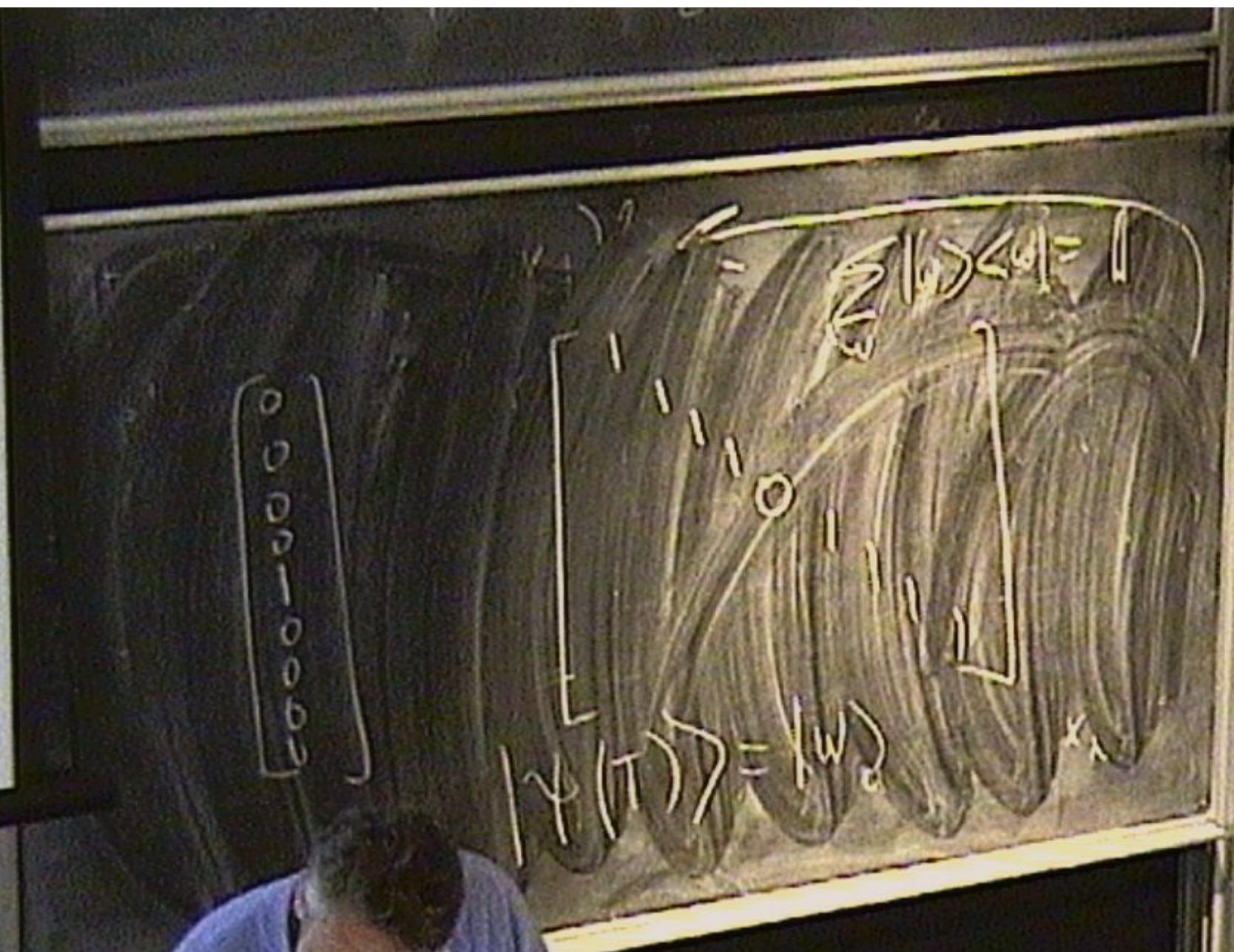
Now $|\psi_\omega, 0\rangle = |\psi_i, 0\rangle$

$$\sum_{\omega} |\langle \psi_\omega, t \rangle - \langle \psi_i, t \rangle|^2 \leq 2E N^{1/2} t$$

for $t = T$

$$N \leq 2E N^{1/2} T$$

$$\frac{N^{1/2}}{2E} \leq T$$



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$$\bullet \frac{d}{dt} |\langle \psi_\omega, t \rangle - \langle \psi_i, t \rangle|^2 = -2 \operatorname{Re} \frac{d}{dt} \langle \psi_\omega, t | \psi_i, t \rangle$$

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$$\bullet \frac{d}{dt} \left| \langle \psi_\omega, t \rangle - |\psi, t\rangle \right|^2 = -2 \Re \frac{d}{dt} \langle \psi_\omega, t | \psi, t \rangle$$

$$= 2 \operatorname{Im} \langle \psi_\omega, t | H_\omega | \psi, t \rangle \\ \leq 2 \left| H_\omega | \psi, t \rangle \right|$$

Now $\sum_3 \epsilon_\omega \langle \psi_\omega \rangle = E$

$$\frac{d}{dt} \sum_3 \left| \langle \psi_\omega, t \rangle - |\psi, t\rangle \right|^2 \leq 2E N^{\gamma_L}$$

$$\bullet \text{Now } |\psi_\omega, 0\rangle = |\psi, 0\rangle$$

$$\sum_3 \left| \langle \psi_\omega, t \rangle - |\psi, t\rangle \right|^2 \leq 2E N^{\gamma_L} t$$

$$\text{for } t = T$$

$$N \leq 2E N^{\gamma_L} T$$
$$\boxed{\frac{N^{\gamma_L}}{2E} \leq T}$$

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$$\bullet \frac{d}{dt} | |\psi_{\omega}, t\rangle - |\psi_i, t\rangle|^2 = -2 \operatorname{Re} \frac{d}{dt} \langle \psi_{\omega}, t | \psi_i, t \rangle$$
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$$\leq 2 | H_{\omega} | \psi_i, t \rangle |$$

Now $\sum_{\omega} : \sum_{\omega} \in |\omega\rangle \langle \omega| = E$

$$\frac{d}{dt} \sum_{\omega} | |\psi_{\omega}, t\rangle - |\psi_i, t\rangle|^2 \leq 2E N^{1/2}$$

$$\bullet \text{Now } |\psi_{\omega}, 0\rangle = |\psi_i, 0\rangle$$

$$\sum_{\omega} | |\psi_{\omega}, t\rangle - |\psi_i, t\rangle|^2 \leq 2E N^{1/2} t$$

$$\text{for } t = T$$

$$N \leq 2E N^{1/2} T$$

$$\boxed{\frac{N^{1/2}}{2E} \leq T}$$

6.

General Search Starting with a One-Dimensional Projector:

$$H_p = \sum_{z \in Z} h(z) |z\rangle$$

$$|s\rangle = \frac{1}{\sqrt{N}} \sum_z |z\rangle$$

$$H_B = E(| - |s\rangle \langle s|)$$

$$|\psi(0)\rangle = |s\rangle$$

Theorem: Let H_p be diagonal in the z -basis with a ground state subspace of dimension K . Let

$$H(t) = (1 - \frac{t}{T}) E(| - |s\rangle \langle s|) + \frac{t}{T} H_p$$

Let P be the projector onto the ground state subspace of H_p and let $b > 0$ be the success probability

$$b = \langle \psi(T) | P | \psi(T) \rangle$$

Then

$$T \geq \frac{b}{E} \sqrt{\frac{N}{K}} - \frac{2\sqrt{b}}{E}$$

7

Simple example

Decoupled problem

$$\text{For each bit } j \quad h_j(z_j) = \begin{cases} 0 & z_j = 0 \\ 1 & z_j = 1 \end{cases}$$

$$h(z) = z_1 + z_2 + \dots + z_n$$

$$\text{Good: } H_B = \sum_j \frac{1}{2} (1 - e^{i\phi_j})$$

~~Ground state of~~ Ground state of H_B is $|s\rangle$

Qubits rotate to winner. $T \sim \sqrt{n}$

$$\text{Bad } H_B = E(|s\rangle\langle s|)$$

T must be greater than $2^{n/2}$

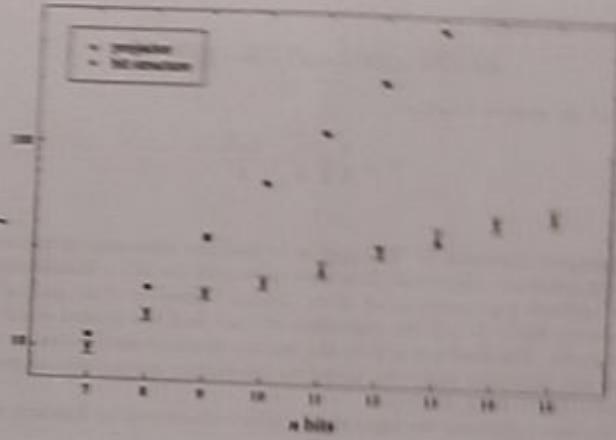


FIG. 1: Median required run time T versus bit number. At each bit number there are 60 random instances of Exact Cover with a single satisfying assignment. We choose the required run time to be the value of T for which quantum adiabatic algorithm has success probability between 0.2 and 0.31. For the projective beginning Hamiltonian we use (8) with $E = n/2$. The plot is log-linear. The error bars show the 95% confidence interval for the true medians.

The total cost function is then

$$h(s) = \sum_i h_i(s).$$

To get H_B to reflect the bit and clause structure we may pick

$$H_{B,i} = \frac{1}{2} \left[(1 - \sigma_x^{(i,j)}) + (1 - \sigma_y^{(i,j)}) + (1 - \sigma_z^{(i,j)}) \right]$$

with

$$H_B = \sum_i H_{B,i}. \quad (14)$$

In this case the ground state of H_B is again $|e\rangle$. With this setup, Theorem 1 does not apply.

We did a numerical study of a particular satisfiability problem, Exact Cover. For this problem if clause c involves bits i_1, i_2 , and i_m , the cost function is

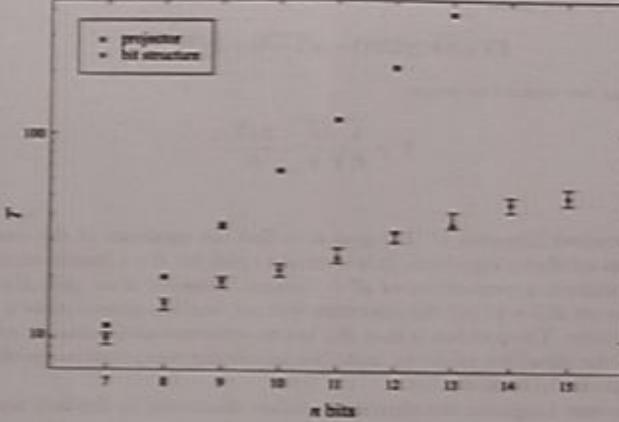


FIG. 1: Median required run time T versus bit number. At each bit number there are 50 random instances of Exact Cover with a single satisfying assignment. We choose the required run time to be the value of T for which quantum adiabatic algorithm has success probability between 0.2 and 0.21. For the projector beginning Hamiltonian we use (8) with $E = n/2$. The plot is log-linear. The error bars show the 95% confidence interval for the true medians.

The total cost function is then

$$h(z) = \sum_c h_c(z).$$

To get H_B to reflect the bit and clause structure we may pick

$$H_{B,c} = \frac{1}{2} \left[(1 - \sigma_x^{(i_c)}) + (1 - \sigma_x^{(j_c)}) + (1 - \sigma_x^{(k_c)}) \right]$$

with

$$H_B = \sum_c H_{B,c}. \quad (13)$$

In this case the ground state of H_B is again $|s\rangle$. With this setup, Theorem 1 does not apply.

We did a numerical study of a particular satisfiability problem, Exact Cover. For this problem if clause c involves bits i_c , j_c and k_c , the cost function is

$$h_c(z) = \int 0 \quad \text{if } z_{i_c} + z_{j_c} + z_{k_c} = 1,$$

7

Simple example

Decoupled problem

$$\text{For each bit } j \quad h_j(z_j) = \begin{cases} 0 & z_j = 0 \\ 1 & z_j = 1 \end{cases}$$

$$h(z) = z_1 + z_2 + \dots + z_n$$

$$\text{Good: } H_B = \sum_j \frac{1}{2} (1 - e^{i\phi_j})$$

~~Ground state of H_B is $|s\rangle$~~

Qubits rotate to winner. $T \sim \sqrt{n}$

$$\text{Bad } H_B = E(|s\rangle \langle s|)$$

T must be greater than $2^{n/2}$

$$i\hbar \frac{d}{dt} |\psi(t)\rangle = H(t) |\psi(t)\rangle$$

$$|\psi(1)\rangle = \left(I - \frac{i\hbar}{\tau} \right)$$

ϵ_{ρ_2}

7

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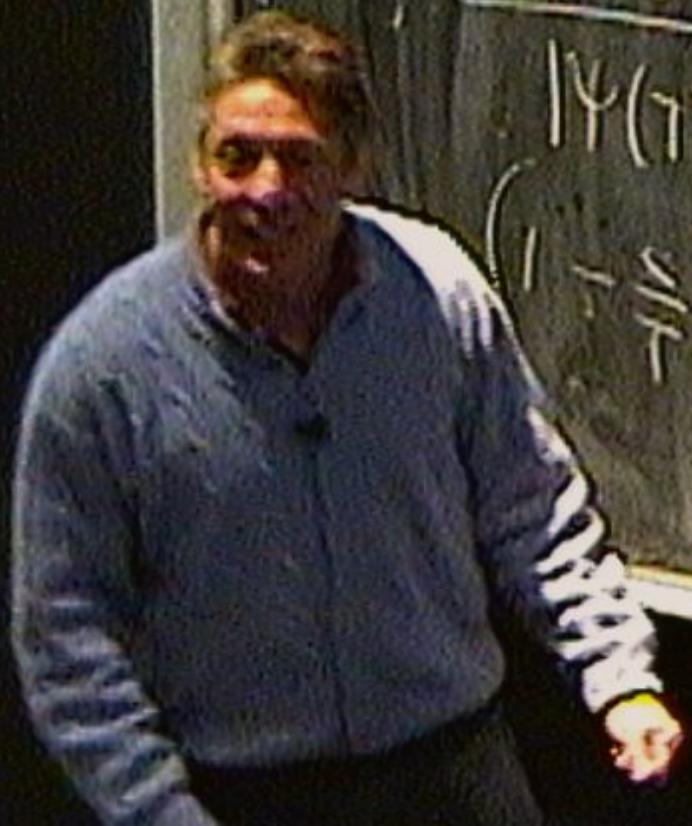
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$$\frac{d|\psi(t)\rangle}{dt} = H(t)|\psi(t)\rangle$$
$$|\psi(1)\rangle = \left(1 - \frac{\pi}{F}\right) |\psi(0)\rangle + \frac{1}{F} |\psi_+\rangle$$



8.

● Search with a Scrambled Problem

Hamiltonian:

Cost function $h(z)$: minimum at $z=0$

Let π be a permutation of $0, 1, \dots, N-1$
(not of the bits!)

$$h^{[\pi]}(z) = h(\pi^{-1}(z)) \text{ . minimum at } \pi(0)$$

$$H_{p,\pi} = \sum_{z=0}^{N-1} h^{[\pi]}(z) |z\rangle\langle z| = \sum_{z=0}^{N-1} h(z) |\pi(z)\rangle\langle\pi(z)|$$

$$\text{Let } H_\pi(t) = H_0(t) + c(t) H_{p,\pi}$$

$$|c(t)| \lesssim 1 \text{ for all } t$$

~~Initial state~~ Evolve from a π
independent state for time T .

$$\therefore \frac{d}{dt} |\psi_\pi(t)\rangle = [H_0(t) + c(t) H_{p,\pi}] |\psi_\pi(t)\rangle$$

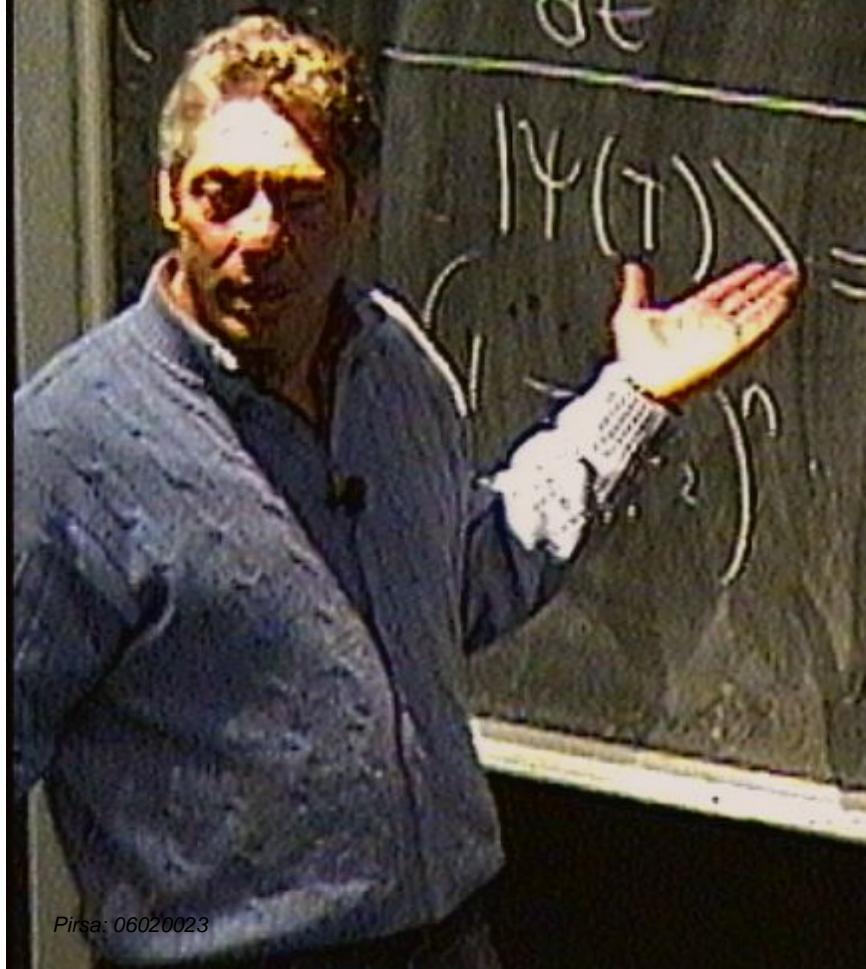
Want $|\psi_\pi(T)\rangle$ near $|\pi(0)\rangle$

$$i\hbar \frac{d}{dt} |\psi(t)\rangle = H(t) |\psi(t)\rangle$$

$$|\psi(t)\rangle =$$

$$\frac{1}{\sqrt{2}}$$

$$\begin{pmatrix} 0 \\ 1 \\ 2 \\ 0 \\ 3 \\ 1 \end{pmatrix}$$



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$$h^{[\pi]}(z) = h(\pi^{-1}(z)) \text{ . minimum at } \pi(0)$$

$$H_{p,\pi} = \sum_{z=0}^{N-1} h^{[\pi]}(z) |z\rangle\langle z| = \sum_{z=0}^{N-1} h(z) |\pi(z)\rangle\langle\pi(z)|$$

$$\text{Let } H_\pi(t) = H_0(t) + c(t) H_{p,\pi}$$

$$|c(t)| \leq 1 \text{ for all } t$$

Evolve from a π
independent state for time T .

$$\therefore \frac{d}{dt} |\psi_\pi(t)\rangle = [H_0(t) + c(t) H_{p,\pi}] |\psi_\pi(t)\rangle$$

Want $|\psi_\pi(T)\rangle$ near $|\pi(0)\rangle$

- Theorem: Suppose a continuous time algorithm of the form on the previous page succeeds with a probability b , $|\langle \hat{T}_\pi(\tau) | \pi(0) \rangle|^2 \geq b$ for a set of $\in N!$ permutations. Then

$$T \geq \frac{\varepsilon^2 b}{16 h^*} \sqrt{N-1} - \frac{\varepsilon \sqrt{\varepsilon/2}}{4 h^*}$$

$$\text{with } h^* = \left(\frac{1}{N-1} \sum_z h^2(z) \right)^{1/2}$$

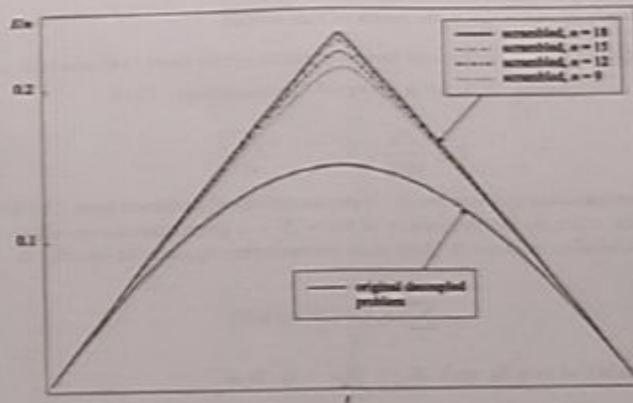


FIG. 2: The scaled ground state energy E/n for a quantum adiabatic algorithm Hamiltonian of a decoupled problem. The lowest curve corresponds to the original decoupled problem. The upper "triangular" curves correspond to single instances of the n -bit decoupled problem, where the problem Hamiltonian was scrambled.

(Since the system is decoupled this is actually the ground state energy of a single qubit.) We then consider the n bit scrambled problem for different values of n . At each n we pick a single random permutation π of $0, \dots, (2^n - 1)$ and apply it to obtain a cost function $h(\pi^{-1}(z))$ while keeping H_B fixed. The ground state energy divided by n is now plotted for $n = 9, 12, 15$ and 18 . From these scrambled problems it is clear that if we let n get large the typical curves will approach a triangle with a discontinuous first derivative at $t = T/2$. For large n , the ground state changes dramatically as t passes through $T/2$. In order to keep the quantum system in the ground state we need to go very slowly near $t = T/2$ and this results in a long required run time.

IV. CONCLUSIONS

In this paper we have two main results about the performance of the quantum adiabatic algorithm when used to find the minimum of a classical cost function $h(z)$ with $z = 0, \dots, N-1$. Theorem 1 says that for any cost function $h(z)$, if the beginning Hamiltonian is a one dimensional projector onto the uniform superposition of all the $|z\rangle$ basis states, the algorithm will not find the minimum of h if T is less than of order \sqrt{N} . This is true regardless of how simple it is to classically find the minimum of $h(z)$.

In Theorem 2 we start with any beginning Hamiltonian and classical cost function h . Replacing $h(z)$ by a scrambled version, i.e. $h^{[\pi]}(z) = h(\pi(z))$ with π a permutation of 0 to $N-1$, will make it impossible for the algorithm to find the minimum of $h^{[\pi]}$ in time less than order \sqrt{N} for a typical permutation π . For example suppose we have a cost function $h(z)$ and have chosen H_B so that the quantum algorithm finds the minimum in time of order $\log N$. Still scrambling the cost function results in algorithmic failure.

These results do not imply anything about the more interesting case where H_B and H_P are

$$G = E$$

$$G = -F$$



$$G = \begin{bmatrix} 2 & 1 \\ 0 & 1 \end{bmatrix}$$

$$G = \begin{bmatrix} 2 & 1 \\ -1 & 1 \end{bmatrix}$$

1
2
3
4
5
6
7
8
9
10



$$G = E^{\beta^{-1}}$$

$$G = -E^{\beta^{-1}}$$

$$\begin{matrix} 0 \\ 1 \\ 2 \\ 3 \end{matrix}$$

$$GH$$

$$H =$$

$$H_{\text{ext}} + SH_F$$

$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

$$H =$$

$$GH \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$$

$$H_F$$

$$GH$$

$$G = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

$$G = -H_F$$

$$H_f(\omega) + S H_F$$

$$G = \begin{bmatrix} E & 0 \\ 0 & 0 \end{bmatrix}$$

$$G = \begin{bmatrix} E & 0 \\ 0 & F \end{bmatrix}$$

$$\begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}$$

$$GH \left\{ \begin{array}{l} H_f \\ GH \end{array} \right.$$

$$H =$$

$$\begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}$$

$$H(s) + sU_f$$

$$G = \begin{bmatrix} 0 & 1 \\ 0 & 0 \end{bmatrix}$$

$$G = -\begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}$$

$$\begin{array}{c} H = \\ \left(\begin{array}{cc} 0 & 1 \\ 1 & 0 \end{array} \right) \\ \text{or} \\ \left(\begin{array}{cc} 0 & 1 \\ 0 & 0 \end{array} \right) \\ \text{or} \\ \left(\begin{array}{cc} 0 & 1 \\ 0 & 0 \end{array} \right) \end{array}$$

GH

H_f

GH_f

$$H(1-s) + s H_F$$

$$G = \left[\begin{array}{c} 1 \\ 0 \\ 0 \\ 0 \end{array} \right]$$

$$G = \left[\begin{array}{c} 1 \\ -1 \\ 0 \\ 0 \end{array} \right]$$

$$H = \left[\begin{array}{c} 0 \\ 0 \\ 0 \\ 0 \end{array} \right]$$

$$GH = \left\{ \left[\begin{array}{c} 0 \\ 0 \\ 0 \\ 0 \end{array} \right] \right\}$$

$$H_f$$

$$GH$$

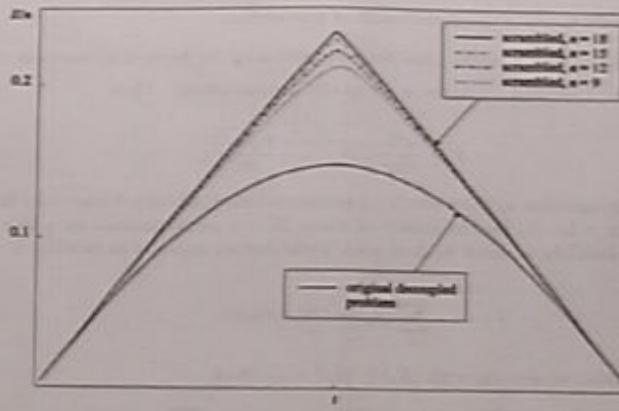


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In this paper we have two main results about the performance of the quantum adiabatic algorithm when used to find the minimum of a classical cost function $h(z)$ with $z = 0, \dots, N - 1$. Theorem 1 says that for any cost function $h(z)$, if the beginning Hamiltonian is a one dimensional projector onto the uniform superposition of all the $|z\rangle$ basis states, the algorithm will not find the minimum of h if T is less than of order \sqrt{N} . This is true regardless of how simple it is to classically find the minimum of $h(z)$.

In Theorem 2 we start with any beginning Hamiltonian and classical cost function h . Replacing $h(z)$ by a scrambled version, i.e. $h^{[\pi]}(z) = h(\pi(z))$ with π a permutation of 0 to $N - 1$, will make it impossible for the algorithm to find the minimum of $h^{[\pi]}$ in time less than order \sqrt{N} for a typical permutation π . For example suppose we have a cost function $h(z)$ and have chosen H_B so that the quantum algorithm finds the minimum in time of order $\log N$. Still scrambling the cost function results in algorithmic failure.

These results do not imply anything about the more interesting case where H_B and H_P are

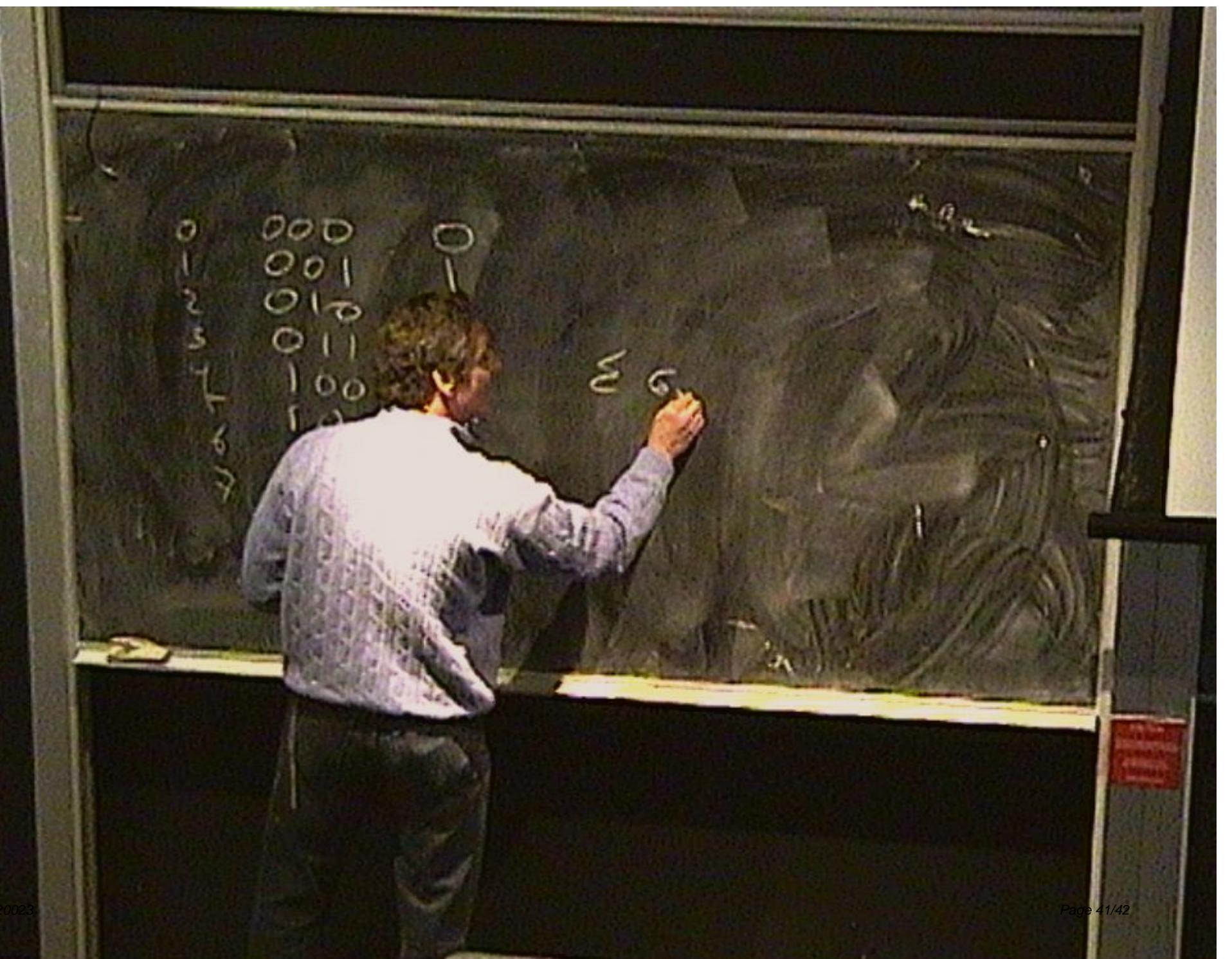
A person wearing a dark long-sleeved shirt is standing in front of a chalkboard, writing binary code with a piece of chalk. The chalkboard has some faint, illegible markings at the top right.

000
001
010
011
100
101

000
001
010
011
100
101
110
111

0
1
2
3
4
5
6
7

= 32



A man in a blue shirt stands at a chalkboard, writing binary numbers. He is positioned on the right side of the frame, facing left. The chalkboard is dark and shows several rows of binary digits written in white chalk. On the left, there are two columns of binary digits: one column has four rows, and the other has three rows. To the right of these columns, there is a single column with five rows. The binary digits are as follows:

Column 1	Column 2	Column 3
0000	000	0
001	001	1
010	010	10
011	011	11
100	100	100
101	101	101
110	110	110
111	111	111

Σ G.