Quantum Timing and Synchronization Problems

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Abstract

Feynman’s model of a quantum computer provides an example of a continuous-time quantum walk. Its clocking mechanism is an excitation of a basically linear chain of spins with occasional controlled jumps which allow for motion on a planar graph. The spreading of the wave packet poses limitations on the probability of ever completing the $s$ elementary steps of a computation: an additional amount of storage space $\delta$ is needed in order to achieve an assigned completion probability. In this note we study the END instruction, viewed as a measurement of the position of the clocking excitation: a $\pi$-pulse indefinitely freezes the contents of the input/output register, with a probability depending only on the ratio $\delta/s$.

Keywords: continuous-time quantum walk; quantum END problem; telomeric chain; $\pi$-pulse trap; quantum subroutines; Grover’s algorithm.

1 Introduction

It has been shown by Feynman[1] that it is possible to implement the sequential application, in the desired order, of a collection $A = A_{s-1} \cdots A_2 A_1$ of unitary operators to an input/output register by using $s$ additional program counter sites.

For the sake of definiteness, we will think of each program counter site $j = 1, 2, \ldots, s$ as occupied by a spin-1/2 system $\tau(j) = (\tau_1(j), \tau_2(j), \tau_3(j))$. We will refer to the collection of such spins, which act in effect as a quantum clocking mechanism, as to a “program line”.

The input/output register will be, similarly, implemented by a collection of a certain number $\mu$ of spin-1/2 systems $\sigma(i) = (\sigma_1(i), \sigma_2(i), \sigma_3(i))$, $i = 1, 2, \ldots, \mu$.

The evolution of the complete system, register + program line, will be given by the Schrödinger equation:

$$i \frac{d}{dt} |\psi(t)\rangle = H |\psi(t)\rangle,$$  \hspace{1cm} (1)

where the Hamiltonian $H$ is supposed to be time independent and involving at most 3-body interactions:

$$H = -\frac{1}{2} \left( \sum_{j=1}^{s-1} \tau_+(j+1) A_j \tau_-(j) + \text{hermitian conjugate} \right).$$  \hspace{1cm} (2)
We have indicated by $\tau_+(j)$ and $\tau_-(j)$ the raising and lowering operators for the third component of the spin occupying the $j$-th program counter site. Since it is requested that each term of the sum is (at most) a 3-body interaction, each $A_j$ either acts on a single spin of the register or is a constant: as shown in Ref.[1], this does not restrict the class of functions computable by the model.

We will restrict our considerations to initial conditions of the form:

$$|\psi(0)\rangle = |\text{register}_0\rangle \otimes |\text{program line}_0\rangle,$$

where $|\text{program line}_0\rangle = |\tau_3(1) = +1, \tau_3(j) = -1 \text{ for } 1 < j \leq s\rangle$ describes the state in which only the spin located at program site number 1 is “up”.

It helps the intuition to think of the initial state $|\text{register}_0\rangle$ of the register as a simultaneous eigenstate of the components of the $\sigma$ spins in selected directions, encoding the initial word (or superposition of words) on which the machine is required to act.

Because of the conservation law $[H, N_3] = 0$, where

$$N_3 = \sum_{j=1}^{s} \frac{1 + \tau_3(j)}{2},$$

the above initial condition has the important consequence that the evolution $|\psi(t)\rangle$ takes place in the $2^s s$ dimensional eigenspace of $N_3$ belonging to the eigenvalue +1.

The intuition of “a single clocking excitation travelling along the program line” emerging from the above conservation law is made precise by introducing the observable position of the excitation, or position of the cursor:

$$Q = \sum_{j=1}^{s} j \frac{1 + \tau_3(j)}{2}.$$

It is then easy to convince oneself that the evolution of the overall system is of the form

$$|\psi(t)\rangle = \sum_{k=1}^{s} c(t, k; s) |\text{register}_{k-1}\rangle \otimes |Q = k\rangle,$$

where, for $1 \leq h \leq s - 1$, $|\text{register}_h\rangle = A_h |\text{register}_{h-1}\rangle$.

In Feynman’s words (adapted to our notations), [3] says that, starting from the initial condition [8], “If at some later time the final site $s$ is found to be in the $|\tau_3(s) = +1\rangle$ state (and therefore all the others in $|\tau_3(j) = -1\rangle$), then the register state has been multiplied by $A_{s-1} \cdots A_2 \cdot A_1$ as desired”. It as been shown in Ref.[2] that this is a somewhat big “If”, under two respects:

i. at no instant of time the probability $|c(t, s; s)|^2$ is larger than $\text{const} \cdot s^{-\frac{3}{4}}$;

ii. the cursor keeps bouncing back and forth between positions 1 and $s$, thus in effect making the above upper bound attainable only at selected instants of time.

The above two statements are reviewed and made quantitative in Section 2. Section 3 is devoted to the “quantum END problem”: we remove the cursor in order to prevent it from returning down the program line and “undoing” the computation. Removing
the cursor and storing the result of the computation in the contents of the register is in effect a measurement procedure, that in Sec. 3 will be modelled by a suitable time dependent perturbation (a $\pi$-pulse) applied to a variant of the Hamiltonian (2). Section 4 is devoted to conclusions and outlook.

2 The Motion of the Cursor

We recall, first of all, that the motion of the cursor does not depend on the operators acting on the register. For the particular case of a sequential program line as the one described by the Hamiltonian (2), this is made evident by the explicit expression of the amplitudes $c(t,k;s)$ in (3): they are given, independently of the operators $A_j$, by:

$$c(t,k;s) = \frac{2}{s + 1} \sum_{n=1}^{s} \exp(it \cos(\theta(n;s))) \sin(\theta(n;s)) \sin(k \theta(n;s)),$$

where:

$$\theta(n;s) = \frac{n\pi}{s + 1}$$

Similar results hold in the case, studied in Ref. [4], in which, because of conditional jumps in the program line (such as the ones needed in the iteration of quantum subroutines), the cursor performs, in effect, a continuous-time quantum walk on a planar graph. In this note we restrict ourselves to the sequential case.

The main purpose of this Section is to give examples of the behaviour recalled in the observations (i) and (ii) of Sec. 1.

This we do with the help of the following Hamiltonian:

$$H = -\frac{1}{2} \left( \sum_{j=1}^{s-1} \tau_+(j+1)A_j \tau_- (j) + \tau_+(s+1)\rho_- \tau_- (s) + \sum_{j=s+1}^{s+\delta-1} \tau_+(j+1) \tau_- (j) + h.c. \right).$$

With respect to the Hamiltonian (2), we have introduced an additional control q-bit $\rho = (\rho_1,\rho_2,\rho_3)$ in the term $\tau_+(s+1)\rho_- \tau_- (s)$; this is an example of a conditional jump in the quantum walk performed by the cursor: it acts non trivially only in the eigenspace belonging to the eigenvalue +1 of $\rho_3$, enabling the transition $|Q = s \rangle \rightarrow |Q = s + 1 \rangle$. If this transition is enabled, then the cursor can visit the additional telomeric sites $s + 1, \ldots, s + \delta$, else it gets reflected back.

Figures 1 and 2 give examples of the behaviour of the probability

$$P_{(s \leq Q)}(t) = P_{(s \leq Q \leq s+\delta)}(t) = \sum_{j=s}^{s+\delta} |\gamma(t,j)|^2$$

of finding the register in the state $A = A_{s-1} \cdots A_2 \cdot A_1 |register_0 \rangle$, under two different initial conditions, which determine two different forms of the amplitudes $\gamma$. 
Figure 1 corresponds to the initial condition $|\text{program line}_0\rangle = |\rho_3 = -1\rangle \otimes |Q = 1\rangle$: the motion of the cursor remains confined to the sites $1, \ldots, s$, as it is $\gamma(t, k) = c(t, k; s)$ if $1 \leq k \leq s$, 0 otherwise. The probability $P_{(s \leq Q)}(t)$ of finding the computation completed satisfies in this case the inequality:

$$ P_{(s \leq Q)}(t) \leq \frac{8}{s^2} \quad (11) $$

Figure 2 corresponds to the initial condition: $|\text{program line}_0\rangle = |\rho_3 = +1\rangle \otimes |Q = 1\rangle$, leading to $\gamma(t, k) = c(t, k; s + \delta)$ for $1 \leq k \leq s + \delta$. For $t$ just below $s + 2\delta$ the probability $P_{(s \leq Q)}(t)$ of finding the computation completed is close to the much less severe upper bound:

$$ P_{(s \leq Q \leq s + \delta)}(t) \leq 1 - \frac{2}{\pi} \left( \arcsin \left( \frac{1}{1 + 2\delta/s} \right) - \left( \frac{1}{1 + 2\delta/s} \right) \sqrt{1 - \left( \frac{1}{1 + 2\delta/s} \right)^2} \right). \quad (12) $$

3 The Quantum END Instruction

The abrupt collapse of $P_{(s \leq Q)}(t)$ at time $t \approx s + 2\delta$, evident from Fig. 2, corresponds to the following fact: travelling with average speed close to 1, at time $t \approx s + 2\delta$ the cursor “returns
down the active part of the program line”, thus, in effect, undoing the calculation.

Bringing the computation to an END, and storing the result is not completely trivial in
the case examined here of a reversible quantum clocking mechanism: “Surely a computer
has eventually to be in interaction with the external world, both for putting data in and for
taking it out![1]”.

A simple model of such interaction is suggested by inspection of the Hamiltonian (9): starting
from the initial condition $|\rho_3 = +1\rangle$, the transition $|Q = s\rangle \to |Q = s + 1\rangle$ is enabled by the
control term $\tau_+ (s + 1) \rho_+ \tau_- (s)$ which, simultaneously, determines the transition $|\rho_3 = +1\rangle \to
|\rho_3 = -1\rangle$.

The transition $|Q = s+1\rangle \to |Q = s\rangle$, enabled by the hermitian conjugate term $\tau_+ (s) \rho_+ \tau_- (s + 1)$,
will be therefore inhibited if, by external means, we enforce the transition $|\rho_3 = -1\rangle \to
|\rho_3 = +1\rangle$ at a time, close to $t_0 = s + 2\delta$, when most of the probability mass is in the region
$s, \ldots, s + \delta$.

Figures 3 (where Figs. 1 and 2 are also reproduced for comparison purpose) presents the
effect of the addition to (9) of the time dependent perturbation

$$h(t) = \frac{1}{2} B(t) \cdot \rho_1$$

(13)

where the “magnetic field” $B(t)$ is non vanishing only in a unit time interval around $t_0$, in
which it takes the value $\pi$. With a probability depending only on the ratio $\delta/s$ (see (12))

Figure 3: The solid line represents the probability of finding the cursor in the telomeric chain
using a $\pi$-pulse applied at time $t_0 = s + 2\delta$. The dashed lines correspond to Figs[1] and 2
between the lengths of the active part and the telomeric part of the program line, the $\pi$-pulse
[13] definitively prevents the cursor from undoing the computation.

The idea of a $\pi$-pulse trap just presented works only if the control q-bit is initialised in the
$|\rho_3 = +1\rangle$ state. It is immediate to convince oneself that the following double trap Hamiltonian
does not suffer from the above limitation:

\[ H = h(t) + \]

\[ -\frac{1}{2} \sum_{j=1}^{s-1} \tau_+(j+1)A_j\tau_-(j) + \sum_{j=s+1}^{s+\delta-1} \tau_+(j+1)\tau_-(j) + \sum_{j=s+\delta+1}^{s+2\delta-1} \tau_+(j+1)\tau_-(j) + \]

\[ + \tau_+(s+1)\tau_-(s) + \tau_+(s+\delta+1)\tau_-(s) + h.c. \]  

(14)

With any initial condition for the control q-bit, under the action of the above Hamiltonian, the \(|\rho_3 = +1\rangle\) component of the state gets definitively trapped in the first telomeric region \(\{s+1, \ldots, s+\delta\}\), the \(|\rho_3 = -1\rangle\) component in the second one \(\{s+\delta+1, \ldots, s+2\delta\}\).

As a final remark of this section, we observe that, acting in effect as a Stern-Gerlach apparatus providing space separation between two different spin states, the term

\[ \text{switch} = (\tau_+(s+1)\tau_-(s) + \tau_+(s+\delta+1)\tau_-(s)) + h.c. \]  

(15)

can be used also to model the preparation ("putting the data in") of a register q-bit in a given spin state.

4 Conclusions and Outlook

Feynman’s time honoured model of a quantum computer and its modern streamlined version, the continuous-time quantum walk\[5\] paradigm (in which the quantum motion of the cursor on a graph is the computation, irrespective of any action on the register), provide a fascinating physical context in which to think of “time” under a quantum perspective and are rich sources of open problems.

Inequality (11) is, for instance, a simple consequence of the spreading of the wave packet in the inertial motion (7) of the cursor on a finite lattice. Is inequality (11) strictly dependent on the model adopted here or is it a hint of a model independent “probability vs. computational complexity” uncertainty relation?

Inequality (12) and the related discussion of Sec. 3 set bounds on the minimum amount of additional space needed in order to have a preassigned probability of storing the result of the complete calculation. Similarly, the preparation of a given input requires time and space in order to perform the required “Stern-Gerlach” preparation. Is it realistic to neglect this costs as it is done in the conventional performance analysis of performance of quantum algorithms?

As a concluding remark, we wish to point out a case, Grover’s algorithm\[6\], in which the nature, classical or quantum, of the clocking agent that successively applies the required primitives does make a difference.

Grover’s algorithm poses the problem of estimating the parameters \(a = (a_1, \ldots, a_\mu) \in \{-1, +1\}^\mu\) appearing in the transformation \(A = 1 - 2P_a\) that an oracle is able to apply to the register. Here we have indicated by \(P_a\) the projector on the state \(|\sigma_3(i) = a_i, \ i = 1, \ldots, \mu\rangle \equiv |a_3\rangle\).

The estimation procedure of \(a\) starts from \(|\text{register}_0\rangle = |\sigma_1(i) = +1, i = 1, \ldots, \mu\rangle \equiv |1_1\rangle\) and proceeds by alternating the action of \(A\) with the action of \(B = 1 - 2|1_1\rangle\langle 1_1|\).

It can be shown\[7\] that:

\[ |\langle a_3| (B \cdot A)^n|1_1\rangle|^2 = \sin^2((2n+1)\theta), \ 	ext{with} \ \theta = \arcsin(2^{-\mu/2}). \]  

(16)
The sharp maximum of (16) at \( n = n_{\text{optimal}} \approx \frac{\pi}{4} 2^{n/2} \) keeps reappearing periodically if the computation proceeds indefinitely after \( n_{\text{optimal}} \) steps.

In Refs. [8] and [9], the oscillatory nature of the overlap (16) between the target state \( |a_3\rangle \) and the current state \((B \cdot A)^n|1_1\rangle\) has been nicely explained in terms of the spectral gap of an analogue Hamiltonian of the form \(-\gamma |a_3\rangle\langle a_3| - |1_1\rangle\langle 1_1|\) acting on the register viewed as an isolated system.

In the context of the model considered in the previous sections, Grover’s algorithm corresponds to the execution of a program line of the form (2) with \( A_j = A \) for odd \( j \), \( A_j = B \) for even \( j \).

Starting from the initial condition \( |\psi(0)\rangle = |1_1\rangle \otimes |Q = 1\rangle \), and taking, for the sake of definiteness, an odd value of \( s \), \( s = 2g + 1 \), it is easy to check that the overlap probability is, in this case, given by

\[
\langle \psi(t)|P_a|\psi(t)\rangle = \sum_{n=0}^{g} \sin^2((2n + 1)\vartheta)(|c(t,2n+1;s)|^2 + |c(t,2n+2;s)|^2) = \\
= \sum_{x=1}^{s} |c(t,x;s)|^2 \sin^2(\vartheta x_{\text{odd}}),
\]

(17)

where \( x_{\text{odd}} \) is the largest odd number not larger than \( x \).

Direct inspection of (17) shows that (16) gives the conditional probability of finding the register in the target state, \( \text{given} \) that the cursor is in the state \( |Q = 2n + 1\rangle \); in loose terms: the oscillatory behaviour (16) describes the overlap probability as a function of the “machine time”, namely the position \( Q \) of the clocking excitation.

![Figure 4: \( \mu = 5; s = 2^\mu + 1 = 65 \).](image)

The right hand side of (17) gives, instead, the overlap probability as a function of the “Galileian time” \( t \), the independent variable in (16).

Figure 4 shows the behaviour of \( \langle \psi(t)|P_a|\psi(t)\rangle \) as a function of \( t \): the coupling of the register with the many clocking degrees of freedom has a damping effect on the oscillation of the state of the register.

In the example just discussed we have used, in fact, a number of clocking degrees of freedom growing exponentially with \( \mu \). It is shown in Refs. [4] and [10] that the above pseudo-dissipative effect can still be observed if, by a careful use of quantum subroutines, the number of program line sites is reduced to a polynomial in \( \mu \).
Acknowledgements

It is a pleasure to thank Professor Alberto Bertoni for his constant attention and encouragement during completion of this work.

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