

2000]Primary 81P68, 93C85; Secondary 94A99.

## Space Searches with a Quantum Robot

Paul Benioff

**ABSTRACT.** Quantum robots are described as mobile quantum computers and ancillary systems that move in and interact with arbitrary environments. Their dynamics is given as tasks which consist of sequences of alternating computation and action phases. A task example is considered in which a quantum robot searches a space region to find the location of a system. The possibility that the search can be more efficient than a classical search is examined by considering use of Grover's Algorithm to process the search results. For reversible searches this is problematic for two reasons. One is the removal of entanglements generated by the search process. The other is that even if the entanglement problem can be avoided, the search process in 2 dimensional space regions is no more efficient than a classical search. However quantum searches of higher dimensional space regions are more efficient than classical searches. Reasons why quantum robots are interesting independent of these results are briefly summarized.

### 1. Introduction

Quantum computers have been the subject of much study, mainly because of computations that can be done more efficiently than on classical computers. Well known examples include Shor's and Grover's algorithms, [1, 2]. Quantum robots have also been recently described as mobile systems, with an on board quantum computer and ancillary systems, that move in and interact with environments of quantum systems [3]. Dynamics of quantum robots are described as tasks consisting of alternating computation and action phases.

Quantum robots can be used to carry out many types of tasks. These range from simple ones such as searching a region of space to determine the unknown location of a system to complex tasks such as carrying out physical experiments. The fact that the spatial searches are similar to the data base searches which are efficiently implemented using Grover's Algorithm [2] suggests that similar results might hold for use of Grover's Algorithm to process results of a quantum search of a spatial region. This is an example of the possible applicability of Grover's algorithm to various physical measurements [6]. If this can be done for the search task, then one would have an example of a task that can be carried out more efficiently by a quantum robot than by any classical robot.

This possibility is analyzed here for a search of a 2 dimensional space region by a quantum robot to locate a system. It is seen that for reversible searches with time independent unitary dynamics, there are two problems preventing the efficient use of Grover's Algorithm. One is that it appears impossible to remove entanglements generated during the search process. The other problem is that the action in the search task, which is the equivalent of the "oracle function", which is assumed in Grover's Algorithm to be evaluated on any argument in one step, takes many

---

<sup>1</sup>1991 *Mathematics Subject Classification.* [.

This work is supported by the U.S. Department of Energy, Nuclear Physics Division, under contract W-31-109-ENG-38.

steps to evaluate. The result is that even if the entanglement problem is ignored, quantum searches of 2 dimensional space regions are no more efficient than classical searches. However quantum searches of higher dimensional space regions are more efficient than classical searches.

The plan of this paper is to give, in the next section, a brief description of quantum robots and a summary of how they are different from quantum computers. This is followed, in Section 2, by a description of the dynamics of tasks as sequences of alternating computation and action phases. An explicit description of the dynamics is given as a Feynman sum over computation-action phase paths. An example of a quantum robot searching a space area to determine the unknown location of a systems is then described, Section 3. The next section is concerned with the use of Grover's algorithm to process the search results. A very brief summary of Grover's Algorithm, in Subsection 4.1, is followed by a description of the problems encountered, Subsection 4.2. The paper finishes with a discussion of why quantum robots are interesting independent of these results.

## 2. Quantum Robots

**2.1. Comparison with Quantum Computers.** Quantum robots are similar to quantum computers in that an important component is an on board quantum computer. Other systems such as a memory system  $m$ , an output system  $o$ , and a control qubit  $c$  are also present in quantum robots [3]. A relatively minor difference is that quantum robots are mobile whereas quantum computers are stationary relative to the environment. For quantum Turing machine models of quantum computers the head moves but the quantum registers are stationary. For networks of quantum gates the qubit systems move but the gates are stationary. This is shown in physical models of interacting qubits. Examples include ion trap models [4] or nuclear magnetic resonance models [5]. In these cases the ion traps and the liquid of active molecules are stationary.

As is the case for quantum computers the effects of the environment on the component systems of the quantum robot need to be minimized. Methods to achieve this include the possible use of shielding or quantum error correction codes [8]. However other than this the dynamical properties of the environment are completely arbitrary.

This is quite different than the case for quantum computers. To see this assume for quantum Turing machines that the quantum registers are the environment of the head. Similarly for gate networks the moving qubits may be considered as the environment of the network of quantum gates. Here the dynamics of these systems is quite restricted in that the states of the registers or moving qubits can change only during interaction with the head or the gate systems. Also the types of changes that can occur are limited to those appropriate for the specific computation being carried out.

This shows the main difference between quantum robots and quantum computers, namely, that for quantum computers the states of the qubits must not change spontaneously in the absence of interactions related to the computation. No such dynamical restrictions apply to the environment of quantum robots. The states of environmental systems may change spontaneously whether the quantum robot is or is not interacting with them.

Another aspect relates to the requirement that the quantum robot cannot be a multistate system with one or a very few number of degrees of freedom, but must include a quantum computer. One reason is the need for computations as part of the implementation of any task (see below). Another is that the quantum robot may need to be able to respond to a large number of different environmental states. Also a large repertoire of different (task dependent) responses to the same environmental state must be available. If the total number  $N$  of needed responses is large then the only physically reasonable approach is to make the number of degrees of freedom of the quantum robot proportional to  $\log N$ . This is satisfied by including a quantum computer on board.

**2.2. Task Dynamics.** The dynamics of quantum robots are described as tasks consisting of alternating computation and action phases [3]. The purpose of each computation phase is to determine what is to be done in the next action phase. The computation may depend on the states of  $o, m$ , and the local environment as input. The goal is to put  $o$  in one of a set  $\mathcal{B}$  of basis states each of which specifies an action. During the computation the quantum robot does not move. Interactions with the environment, if any, are limited to local entanglement interactions of the type occurring in the measurement process (premeasurements in the sense of Peres [7]).

The purpose of an action phase is to carry out the action determined by the previous computation phase. The action is determined by the state of  $o$  and may include local premeasurements of the environment state. Activities during this phase include motion of the quantum robot and local changes in the state of the environment. The action is independent of the state of the on board quantum computer and the  $m$  system. If  $o$  is in a state in  $\mathcal{B}$ , then the state does not change during the action.

The purpose of the control qubit  $c$  is to regulate which type of phase is active. The computation [action] phase is active only if  $c$  is in state  $|1\rangle$ ,  $[|0\rangle]$ . Thus the last step of the computation [action] phase is the change  $|1\rangle \rightarrow |0\rangle$ ,  $[|0\rangle \rightarrow |1\rangle]$ .

The overall system evolution is described here using a discrete space time lattice. In this case a unitary elementary step operator  $\Gamma$  gives the overall system evolution during an elementary time step  $\Delta$ . Since  $\Gamma$  has, in general, nonzero matrix elements between environmental degrees of freedom and quantum robot degrees of freedom, it describes the evolution of the environment and the quantum robot as well as interactions between the environment and the quantum robot.

It is useful to decompose  $\Gamma$  into two terms based on the states of the control qubit. If  $P_0^c$  and  $P_1^c$  are projection operators on the respective control qubit states  $|0\rangle$  and  $|1\rangle$ , then

$$(2.1) \quad \Gamma = \Gamma(P_0^c + P_1^c) = \Gamma_a + \Gamma_c.$$

Here  $\Gamma_a$  and  $\Gamma_c$  are step operators for the action and computation phases. Interactions among environmental degrees of freedom as well as degrees of freedom of the quantum robot other than those taking part in the task dynamics, if any, are also included in both operators.

Some of the conditions described for the computation and action phases are reflected in properties that the operators  $\Gamma_a$  and  $\Gamma_c$  must satisfy. In particular, if  $P_x^{QR}$ ,  $P_d^o$  are projection operators for finding the quantum robot at each lattice position  $x$  ( $x = x_1, x_2, \dots, x_d$  in  $d$ -dimensional space) and the output system in

any state  $|d\rangle$  in  $\mathcal{B}$ , then one has

$$(2.2) \quad \begin{aligned} \Gamma_c P_x^{QR} &= P_x^{QR} \Gamma_c \\ \Gamma_a P_d^o &= P_d^o \Gamma_a \end{aligned}$$

These commutation relations express the requirements that the position of the quantum robot does not change during the computation phase and, except for possible entanglements with environmental states, the state of the output system is not changed during the action phase. These entanglements would occur if  $o$  was in a linear superposition of  $\mathcal{B}$  states each of which resulted in different environment states during the action. Another property of  $\Gamma_a$  is that it is the identity operator on the space of states for the on board quantum computer and memory system degrees of freedom. Note also that  $\Gamma_a$  and  $\Gamma_c$  do not commute.

If  $\Psi_0$  is the overall system state at time 0 then the state at time  $n\Delta$  is given by  $\Psi_n = (\Gamma_a + \Gamma_c)^n \Psi_0$ . The amplitude for finding the quantum robot and environment in a state  $|w', j\rangle$  is given by

$$(2.3) \quad \Psi_n(w', j) = \sum_{w, i} \langle w', j | (\Gamma_a + \Gamma_c)^n | w, i \rangle \Psi_0(w, i).$$

Here  $|w\rangle, |w'\rangle$  denote the states of all environmental and quantum robot systems except the control qubit in some suitable basis, and  $i, j = 0, 1$  refer to the states of  $c$ .

All the information about the dynamics of the system is given in the matrix elements  $\langle w', j | (\Gamma_a + \Gamma_c)^n | w, i \rangle$ . For each  $w, w', n, i, j$  the matrix element can be expanded in a Feynman sum over phase paths [3, 9]. One first expands  $(\Gamma_a + \Gamma_c)^n$  as a sum of products of  $\Gamma_a$  and  $\Gamma_c$ :

$$(2.4) \quad \begin{aligned} (\Gamma_a + \Gamma_c)^n &= \sum_{v_1=a,c} \sum_{t=1}^n \sum_{h_1, h_2, \dots, h_t=1}^{\delta(\sum, n)} (P_0^c + P_1^c) (\Gamma_{v_t})^{h_t} (\Gamma_{v_{t-1}})^{h_{t-1}}, \\ &\dots, (\Gamma_{v_2})^{h_2} (\Gamma_{v_1})^{h_1}. \end{aligned}$$

In this expansion the number of phases is given by the value of  $t$  which ranges from  $t = 1$  corresponding to one phase with  $n$  steps to  $t = n$  corresponding to  $n$  alternating phases each of 1 step. The duration of the  $\ell$ th phase is given by the value of  $h_\ell$  for  $\ell = 1, 2, \dots, t$ . The requirement that the total number of steps equals  $n$ , or  $h_1 + h_2 + \dots + h_t = n$ , is indicated by the upper limit  $\delta(\sum, n)$  on the  $h$  sum. The alternation of phases is shown by  $v$  where  $v_{m+1} = a$  (or  $c$ ) if  $v_m = c$  (or  $a$ ). The factor  $P_0^c + P_1^c$  expresses the fact that the  $t$ th phase may not be completed.

Expansion in a complete set of states between each of the phase operators  $(\Gamma_{v_\ell})^{h_\ell}$  gives the desired path sum:

$$(2.5) \quad \begin{aligned} \langle w', j | (\Gamma_a + \Gamma_c)^n | w, i \rangle &= \sum_{t=1}^n \sum_{p_2, \dots, p_t} \sum_{h_1, h_2, \dots, h_t=1}^{\delta(\sum, n)} \langle w', j | (\Gamma_{v_t})^{h_t} | p_t \rangle \\ &\dots, \langle p_3 | (\Gamma_{v(2)})^{h_2} | p_2 \rangle \langle p_2 | (\Gamma_i)^{h_1} | w, i \rangle \end{aligned}$$

Here the sum is over all paths  $p$  of states of length  $t+1$  with beginning and endpoints given by the states  $|w\rangle$  and  $|w'\rangle$ . That is  $|p_1\rangle = |w\rangle, |p_t\rangle = |w'\rangle$ . The states of the control qubit have been suppressed as they correspond to the values of  $v$ . Note that  $v(1) = i$ .

FIGURE 1. A Schematic Model of a Quantum Robot at the Origin of  $R$ . Both the memory system lattice (m) and Quantum Turing machine lattice  $\mathcal{L}$  are shown with the head  $h$  that moves on the lattices. The control qubit (c) and output system (o) are also shown. The quantum robot is greatly magnified relative to  $R$  to show details.

Each term in this large sum gives the amplitude for finding  $t$  alternating phases in the first  $n$  steps where the  $\ell$ th phase begins with all systems (except for c) in state  $|p_\ell\rangle$  and ends after  $h_\ell$  steps in state  $|w_{\ell+1}\rangle$ . The sums express the dispersion in the duration or number of steps in each phase ( $h$  sums), in the number of phases ( $t$  sum), and in the initial and terminal states for each phase ( $p$  sums).

### 3. An Example of Quantum Searching

Quantum robots are well suited for carrying out search tasks. As a simple example consider a search task where a quantum robot searches a large square area  $R$  of  $N \times N$  sites to locate a system  $s$ . To keep things simple  $s$  is assumed to be motionless and located at just one unknown site. The goal of the search is to determine the location of  $s$  in  $R$ .

The quantum robot consists of an on board quantum computer, memory and output systems, and a control qubit. The on board computer is assumed here to be a quantum Turing machine consisting of a head moving on a cyclic lattice of  $O(\log N)$  qubits.  $O(-)$  denotes of the order of. The memory system also is a cyclic lattice which is taken here to have about the same number of qubits and to lie adjacent to the computation lattice. A schematic representation of the quantum robot located at a corner (the origin) of  $R$  is shown in the figure.

One method of carrying out the search is to let the coordinates  $X, Y$  with  $0 \leq X, Y \leq N-1$  of each point of  $R$  define a search path. If the memory is initially in state  $|X, Y\rangle_m$  the quantum robot, starting from the location  $0, 0$  moves  $X$  sites

in the  $x$  direction, then  $Y$  sites in the  $y$  direction and looks for  $s$  at its location. After recording the presence or absence of  $s$  at the site and further processing, if any, the quantum robot returns along the path to the origin.

A more detailed description starts with the qubits in the memory,  $m$ , and computation lattice,  $\mathcal{L}$ , in the state  $|X, Y\rangle_m |0\rangle_{\mathcal{L}}$  the output system  $o$  in state  $|dn\rangle_o$  and a computation phase active ( $c$  in state  $|1\rangle_c$ ). After copying the  $m$  state onto  $\mathcal{L}$  to give the state  $|X, Y\rangle_m |X, Y\rangle_{\mathcal{L}}$ , the computation phase checks to see if  $X = 0$  or  $X > 0$ . If  $X > 0$  the computation phase continues by subtracting 1 from  $|X, Y\rangle_{\mathcal{L}}$  to give  $|X - 1, Y\rangle_{\mathcal{L}}$ . It ends by changing the  $o$  state to  $|+x\rangle_o$  and the  $c$  state to  $|0\rangle_c$ .

The action phase consists of one step (one iteration of  $\Gamma_a$ ) in which the quantum robot moves one lattice site in the  $+x$  direction and the  $c$  state is converted back to  $|1\rangle_c$ . The process is repeated until the state with  $X = 0$  is reached on  $\mathcal{L}$ . Then the above process is repeated for  $Y$  (the  $o$  state now becomes  $|+y\rangle_o$  to denote one step motion in the  $+y$  direction) until  $Y = 0$  is reached in the state of  $\mathcal{L}$ .

At this point the presence or absence of  $s$  at the location  $X, Y$  of the quantum robot is recorded during a computation phase and, after further processing, if any, the quantum robot returns along the same path. This is done by interleaving motion of the quantum robot in the  $-y$  and  $-x$  directions, with corresponding  $o$  states  $|-y\rangle_o, |-x\rangle_o$ , with adding 1 to the  $y$ , and then  $x$  components of the  $\mathcal{L}$  state with checking if the values  $Y$  and then  $X$  are reached. This is done by stepwise comparison with the state of  $m$  which remains unchanged.

When the state of  $\mathcal{L}$  is the same as the state of  $m$  the quantum robot has returned to the starting point at the origin of  $R$ . A computation phase changes the  $o$  state to  $|dn\rangle_o$  and transfers motion to some ballast system. As has been noted this is necessary to preserve reversibility and the corresponding unitarity of the dynamics [10].

Examples of ballast motion consist of repetitions of adding 1 to a large lattice of  $M$  qubits or emitting a particle which moves away from  $R$ . In the first case with a finite number  $2^M$  of ballast states, the quantum robot remains in the final state of the search degrees of freedom for a finite time only before the search process is undone. This does not occur for the second case with an infinite number of ballast states.

#### 4. Grover's Algorithm and the Quantum Search

Before applying Grover's Algorithm to process the results of the search, it is useful to understand what it does and how it works. A very brief summary, that follows Grover [2] and Chen et al [11], is given next.

**4.1. Grover's Algorithm.** Suppose one has a data base  $B$  of  $N$  elements and a function  $f$  that takes the value 0 on all elements except one,  $\omega$ , on which  $f$  has value 1. It is assumed that  $\omega$  is completely unknown and that a procedure is available for obtaining the value of  $f$  on any element of the data base in 1 step. Let each  $x$  in  $B$  correspond to a unique length  $n$  binary string and  $|x\rangle_B$  be the corresponding  $n$ -qubit state.

Let the initial state for the search be given by

$$(4.1) \quad \phi = \frac{1}{\sqrt{N}} \sum_{x \in B} |x\rangle$$

where the sum is over all  $N$  elements  $x$  in  $B$ . This corresponds to a coherent sum over all product  $|0\rangle, |1\rangle$  states of  $n$  qubits in a quantum computer if  $N = 2^n$ . This state is easily constructed from the constant 0 state  $|\underline{0}\rangle = \otimes_{j=1}^n |0\rangle_j$  by applying the operator  $(1/\sqrt{2})(\sigma_z + \sigma_x)$  to each qubit. Here  $\sigma_x, \sigma_z$  are the Pauli matrices. This is referred to as the Walsh-Hadamard transformation  $W$ . Thus  $\phi = W|\underline{0}\rangle$ .

Define the unitary operator  $Q$  by  $Q = -I_\phi I_\omega$  where  $I_\phi = 1 - 2P_\phi$  and  $I_\omega = 1 - 2P_\omega$ . Both  $P_\phi$  and  $P_\omega$  are projection operators on the states  $\phi$  and  $|\omega\rangle$ . Let  $|\alpha\rangle = (1/\sqrt{N-1}) \sum_{x \neq \omega} |x\rangle$  be the coherent sum over all states  $|x\rangle$  with  $x$  in  $B$  and different from  $\omega$ . Since  $|\alpha\rangle$  and  $|\omega\rangle$  are orthonormal they form a binary basis for a 2 dimensional Hilbert space.

One can expand  $\phi$  in this basis:

$$(4.2) \quad \phi = \sqrt{\frac{N-1}{N}} |\alpha\rangle + \frac{1}{\sqrt{N}} |\omega\rangle.$$

In the same basis  $Q$  has the representation

$$(4.3) \quad Q = \begin{pmatrix} \cos \theta & \sin \theta \\ -\sin \theta & \cos \theta \end{pmatrix}$$

where  $\cos \theta = 1 - 2/N$  and  $\sin \theta = 2\sqrt{N-1}/N$ .

This shows that  $Q$  acting on  $\phi$  corresponds to a rotation by  $\theta$ , and  $m$  iterations of  $Q$  correspond to a rotation by  $m\theta$ . So, carrying out  $m$  iterations of  $Q$  on  $\phi$  where  $m\theta \approx \pi/2$ , rotates  $\phi$  from a state that is almost orthogonal to  $|\omega\rangle$  to a state that is almost parallel to  $|\omega\rangle$ . Measurement of this final state gives with high probability, the value of  $\omega$ .

Grover's algorithm derives its efficiency from the fact that this rotation is achieved with  $m \sim \sqrt{N}$  whereas classically  $\sim N$  steps are needed to find  $\omega$  with high probability. The iteration of  $Q$  must be stopped at the right value of  $m$  because additional iterations will continue to rotate  $\phi$ .

Efficient implementation of this algorithm on a quantum computer, corresponds to iteration of  $Q$  on each component of  $\phi$ . This requires that it is possible to determine, in a small number of steps, if  $x = \omega$  or  $x \neq \omega$ . This is often described in terms of an unknown or "oracle" function  $f$  on  $B$ , where  $f(x) = 0 [= 1]$  if  $x \neq \omega$  [ $x = \omega$ ], that can be evaluated in one step on any  $x$  in  $B$ .  $I_\phi$  is implementable in  $O(n)$  steps as  $\phi = W|\underline{0}\rangle$  and  $W$  is the product of  $n$  single qubit operators.

Grover [2, 6] first introduced the algorithm for searching an unstructured data base of  $N = 2^n$  elements for a single element (f has value 1 on just one element). Since then the algorithm has been much studied under various generalizations. These include searches for several elements (f has value 1 on several elements) [11, 12, 13], searches in which  $N$  is arbitrary [12], and searches in which the initial amplitude distribution of the component states is arbitrary [14]. It has also been shown that the algorithm is optimally efficient [15]. Further development is described in other work [16, 17]. However, as has been recently emphasized [18], all these searches depend on the fact that the evolving state is and remains a coherent superposition of components corresponding to elements of the data base.

**4.2. Problems with the Use of Grover's Algorithm.** The description in Section 3 of the quantum search was for the initial memory state  $|X, Y\rangle_m$ . However

if the memory system lattice is in the initial state

$$(4.4) \quad \psi_m = (1/N) \sum_{X,Y=0}^{N-1} |X,Y\rangle_m,$$

then the description also applies to each component state  $|X,Y\rangle$ .

As was the case for the initial state  $\phi$ , the state  $\psi_m$  can be efficiently prepared from the state  $|0\rangle_m$  using the Walsh-Hadamard transformation.  $(1/\sqrt{2})(\sigma_z + \sigma_x)$  on each qubit of  $m$ . For this initial memory state all  $N^2$  searches are carried out coherently. Since the path lengths range from 0 to  $2N$ , the quantum robot can search all sites of  $R$  and return to the origin in  $O(N \log N)$  steps. Since this is less than the number of steps,  $O(N^2 \log N)$ , required by a classical robot, the question arises if Grover's algorithm can be used to process the final memory state to determine the location of  $s$ . If this is possible, the overall search and processing should require  $O(N \log N)$  steps which is less than that required by a classical robot.

It is worth a digression at this point to see that Grover's Algorithm is not applicable to the usual method of recording the presence or absence of  $s$  at a site. To see this assume that  $s$  is at site  $X_0, Y_0$  and that an extra qubit,  $r$  of the memory is set aside to record the presence or absence of  $s$ . If  $r$  is initially in state  $|0\rangle_r$  and is changed to state  $|1\rangle_r$  only in the presence of  $s$ , then the initial memory state  $\phi_I = (1/N) \sum_{X,Y=0}^{N-1} |X,Y\rangle_m |0\rangle_r$  is changed to the final memory state

$$(4.5) \quad \phi_f = (1/N) \left( \sum_{X,Y \neq X_0 Y_0} |X,Y\rangle_m |0\rangle_r + |X_0, Y_0\rangle_m |1\rangle_r \right).$$

after the quantum robot has returned to the origin of  $R$ .

The idea then would be to use Grover's algorithm [2] by carrying out  $N$  iterations of a unitary operator  $U$  to amplify the component state  $|X_0, Y_0\rangle_m |1\rangle_r$  at the expense of the other components. Following Grover and others [2, 11], define  $U$  by  $U = -I_{\phi_i} I_{|1\rangle_r}$  where  $I_{\phi_i} = 1 - 2P_{\phi_i}$  and  $I_{|1\rangle_r} = 1 - 2P_{|1\rangle_r}$ . Here  $P_{\phi_i}$  and  $P_{|1\rangle_r}$  are projection operators on the memory state  $\phi_i$  and the record state  $|1\rangle_r$ .  $P_{|1\rangle_r}$  is the identity on other memory degrees of freedom.

It is clear that for this case iterations of  $U$  can be carried out efficiently. However, the problem is that the initial state  $\phi_i$  contains a component,  $|X_0, Y_0\rangle_m |0\rangle_r$ , not present in the final state  $\phi_f$ , Eq. 4.5. Also the state  $\phi_i$  does not contain the state  $|X_0, Y_0\rangle_m |1\rangle_r$ . In this case  $U$  does not have a two dimensional representation in the basis pair

$$(4.6) \quad \frac{1}{\sqrt{N^2 - 1}} \sum_{X,Y \neq X_0 Y_0} |X,Y\rangle_m |0\rangle_r ; \quad |X_0 Y_0\rangle_m |1\rangle_r$$

obtained from Eq. 4.5. As a result  $U$  cannot be represented as a rotation that, under iteration, rotates the desired component to be almost parallel to the initial state.

This problem can be avoided by changing the method of recording the presence or absence of  $s$  and not using an extra qubit  $r$ . Here, following Grover [2, 6], the sign of the component corresponding to the location of  $s$  is changed. In this case the initial memory state, Eq. 4.4, becomes the final memory state

$$(4.7) \quad \psi_f = (1/N) \left( \sum_{X,Y \neq X_0 Y_0} |X,Y\rangle_m - |X_0, Y_0\rangle_m \right).$$



after return of the quantum robot. In this case  $U$  is defined as  $U = -I_{\psi_m} I_{X_0 Y_0}$  where  $I_{\psi_m} = 1 - 2P_m$  and  $I_{X_0 Y_0} = 1 - 2P_{X_0 Y_0}$ . Here  $P_m$  and  $P_{X_0 Y_0}$  are projection operators on the memory states  $\psi_m$  and  $|X_0 Y_0\rangle_m$ .

The problem here is that the only way to determine the action of  $I_{X_0 Y_0}$  is by repeating the search part of the process. This is not efficient as it requires  $O(N \log N)$  steps. In the language of much of the work on Grover's algorithm this corresponds to the fact that it requires  $O(N \log N)$  steps to determine the value of the oracle function instead of just one step as is usually assumed. In this case the advantage of quantum over classical searching is lost for 2 dimensional regions as use of Grover's Algorithm would require  $O(N)$  searches each requiring  $O(N \log N)$  steps.

This suggests that a method be considered in which the Grover iterations are done prior to return when the quantum robot is at the path endpoint. At this point the component memory states are entangled with the quantum robot position states as the overall state has the form  $(1/N) \sum_{X,Y} |X, Y\rangle_m |X, Y\rangle_{QR}$  where  $|X, Y\rangle_{QR}$  is the quantum robot position state for the site  $X, Y$ .

In this case  $I_{X_0 Y_0}$  can be efficiently carried out on each initial component memory state  $|X, Y\rangle$  by a local observation to see if  $s$  is or is not at the site  $X, Y$ . Also the action of  $U = -I_{\psi_m} I_{X_0 Y_0}$  on each component memory state is given by

$$(4.8) \quad U|X, Y\rangle_m = -I_m |X, Y\rangle_m = \frac{2}{\sqrt{N}} \psi_m - |X, Y\rangle_m$$

if  $|X, Y\rangle_m \neq |X_0, Y_0\rangle_m$  and

$$(4.9) \quad U|X_0, Y_0\rangle_m = I_m |X_0, Y_0\rangle_m = -\frac{2}{\sqrt{N}} \psi_m + |X_0, Y_0\rangle_m$$

if  $|X, Y\rangle_m = |X_0, Y_0\rangle_m$ . Here, as before,  $\psi_m = (1/N) \sum_{X,Y=0}^{N-1} |X, Y\rangle_m$ .

Here the problem is that there is no efficient way to carry out more than one iteration of  $U$ . As noted above the first iteration can be done efficiently. However additional iterations require that the action of  $I_{X_0 Y_0}$  on memory component states  $|X', Y'\rangle$  be evaluated for arbitrary values of  $X', Y'$  while the quantum robot remains at site  $X, Y$ . This cannot be done efficiently as the quantum robot has no way of knowing whether  $s$  is or is not at these different locations. To know this the quantum robot must go to the site  $X', Y'$  to see if  $s$  is there. This is inefficient as such a trip requires  $O(N \log N)$  steps. (Actions are efficient if they require  $O(\log N)$  steps or less. Low powers of  $\log N$  are also acceptable.)

One sees from this that implementation of Grover's Algorithm using either of these methods requires  $O(N)$  iterations of  $U$  (as  $R$  has  $N^2$  sites) where each iteration requires  $O(N \log N)$  steps. The resulting number of steps required,  $O(N^2 \log N)$ , is the same as that needed by a classical robot. So quantum searches of 2 dimensional space regions combined with Grover's Algorithm are no more efficient than classical searches.

It is of interest to note that quantum searches of higher dimensional space regions combined with Grover's Algorithm are more efficient than classical searches. To see this assume a search of a  $d$  dimensional cube of  $N^d$  sites with the memory in the initial state

$$(4.10) \quad \Psi = \frac{1}{N^{d/2}} \sum_{X_1, \dots, X_d=0}^{N-1} |X_1, X_2, \dots, X_d\rangle_m$$

Carrying out Grover's Algorithm requires  $O(N^{d/2})$  iterations of  $U$  where, as before, each iteration of  $U$  requires  $O(N \log N)$  steps. This follows from the fact that the number of dimensions appears as a multiplicative factor for the number of steps. Also  $O(dN \log N) = O(N \log N)$ . So the overall process requires  $O(N^{(d/2)+1} \log N)$  steps. For  $d > 2$  this is more efficient than a classical search requiring  $O(N^d \log N)$  steps.

The discussion so far has ignored the entanglement problems. These problems, which apply to all the above cases, result from the fact that the task evolution, starting from the initial unentangled product state  $\psi_m|0\rangle_{\mathcal{L}}|0,0\rangle_{QR} \cdots$ , generates entanglements between the position states  $|X', Y'\rangle_{QR}$  of the quantum robot and the components  $|X, Y\rangle_m$  of the memory state  $\psi_m$ . In order for Grover's Algorithm to work it is necessary to remove this entanglement at the end of each search cycle or iteration of  $U$  so that the final memory state is  $\psi_f$ <sup>1</sup>.

This entanglement occurs because the unitary dynamics is reversible and the number of steps needed to complete the search task is different for different component states of  $\psi_m$ . Here the number ranges from  $O(1)$  for the path  $|0,0\rangle_m$  to  $O(2N)$  for the path  $|N-1, N-1\rangle_m$ . This means that the various components of the quantum robot complete an iteration of  $U$  at different times. This is independent of whether the Algorithm is completed after or prior to return.

Because of the reversibility each component cannot simply stop and wait until the longest search component is complete. It must instead embark on motion of irrelevant or ballistic degrees of freedom. This means that the memory components  $|X, Y\rangle_m$  exchange entanglement with the quantum robot position states for entanglement with states of ballistic degrees of freedom. Since use of Grover's Algorithm requires the removal of this entanglement [18], the question arises whether it is possible to insert delays into each of the memory components that are computed, for example, after the quantum robot returns to the origin of  $R$  at the end of each cycle. If this works then there would be some time or step number at which the entanglement is removed and the original product structure of the initial state recovered, with  $\psi_f$  replacing  $\psi_m$ .

This use of delays to remove the entanglements reversibly requires that no memory of the magnitude of the delay be left in the delay degrees of freedom. Otherwise one ends up with entanglement with the delay degrees of freedom. Also determining the magnitude of each delay is not trivial as it depends not only on the lengths of each of the paths but on the number of steps in the computation phases used to determine motion along the paths. This includes the dependence of the number of steps required to subtract 1 from a number  $M$  on the value of  $M$  (through the number of "carry 1" operations needed [20]).

Based on these considerations it seems doubtful that one can use Grover's Algorithm to efficiently process the results of a quantum search of a space region  $R$ . Even if the entanglement problem were solvable, the above results show that, for 2 dimensional space regions, use of Grover's Algorithm is no more efficient than a classical search. For higher dimensional searches the Algorithm is more efficient. Note that this conclusion is independent of the details of the quantum robot. It applies to any quantum system such as a mobile head that contains

---

<sup>1</sup>The entanglement referred to here is different from that of the memory state qubits. The latter is generated during iteration of the Grover operator and is necessary for successful operation of Grover's Algorithm on multiqubit states [19].

sufficient information on board to tell it where to go, what to do on arrival at the endpoint, and how to return to the origin.

## 5. Discussion

In spite of these pessimistic results, quantum robots are interesting objects of study. For instance they may be useful test beds for study of control of quantum systems [21, 22] as the dependence of the task dynamics on the local environmental state is, for some tasks, similar to a feedback loop.

Quantum robots and the associated task dynamics also make clear what is and is not being done in any task. This is shown by the quantum search task in that the quantum robot does no monitoring or control of its behavior. It (or the on board quantum computer) has no knowledge of where it is in  $R$  at any point or even if it is in  $R$ . For each component memory state  $|X, Y\rangle_m$  there are  $X$  computation phases with the output system  $o$  in state  $|+x\rangle_o$  and  $Y$  phases with  $o$  in state  $|+y\rangle_o$ . These phases are interspersed with  $X$  and  $Y$  action phases during which anything can happen. For example the quantum robot might move outside  $R$  or it might not move at all. Of course for these cases it is unlikely that the quantum robot would return to the origin at the end of the task.

This illustrates a valuable aspect of the description of the task dynamics of quantum robots as sequences of alternating computation and action phases. This is that, for the search task examples described here, it makes very clear the lack of awareness and control the quantum robot has over what has happened in the action phases and what it is doing. This argument applies to the computation phases also. For example the "subtract 1" steps could carry out an arbitrary change to the memory state and the task would continue. In this case the task would no longer be a search task but would be something else.

These considerations are also part of foundational reasons why quantum robots and quantum computers are interesting. If quantum mechanics (or some extension such as quantum field theory) is assumed to be universally applicable, then all systems involved in the validation of quantum mechanics are quantum systems. This includes the systems that make theoretical computations (which includes quantum computers) and the systems that carry out experiments (which includes quantum robots). Thus, in some sense quantum mechanics must describe its own validation, to the maximum extent possible. Exploration of this and the questions of self consistency and possible incompleteness that may occur make this an interesting path of inquiry.

In addition quantum robots, and to some extent quantum computers, are natural systems for investigating several questions. In particular what physical properties must a quantum system have such that

- It is aware of its environment?
- It has significant characteristics of intelligence?
- It changes states of some quantum systems so that the new states can be interpreted as text having meaning to the system generating the text [23]?

In addition there is a sense in which the existence problem for quantum systems having all these properties is already solved. That is, these systems include the readers, and hopefully the author, of this paper.

## References

- [1] P. Shor, in *Proceedings, 35th Annual Symposium on the Foundations of Computer Science*, S. Goldwasser (Ed), IEEE Computer Society Press, Los Alamitos, CA, 1994, pp 124-134; SIAM J. Computing, **26** (1997), 1481-1509.
- [2] L.K.Grover, in *Proceedings of 28th Annual ACM Symposium on Theory of Computing* ACM Press New York 1996, pp. 212-219; Phys. Rev. Letters, **79** (1997), 325-328; G. Brassard, Science **275** (1997), 627-628.
- [3] P. Benioff, Phys. Rev. A **58** (1998), 893-904; *Quantum Robots* in, Feynman and Computation, Exploring the Limits of Computers, Anthony Hey, Ed, Perseus Books, Reading, MA. 1998; Los Alamos Archives Preprint quant-ph/9807032.
- [4] J. I. Cirac and P. Zoller, Phys. Rev. Letters **74** (1995), 4091-4094; C. Monroe, D. M. Meekhof, B. E. King, W. M. Itano, and D. J. Wineland, Phys Rev. Letters **75** (1995), 4714-4717; P. Domokos, J. M. Raimond, M Brune, and S. Haroche, Phys. Rev. A **52** (1995), 3554-3559; C. Monroe, D. Leibfried, B. E. King, D. M. Meekhof, W. M. Itano, and D. J. Wineland, Phys. Rev. A **55** (1997), R2489-2491; J. F. Poyatos, J. I. Cirac, and P. Zoller, Phys. Rev. Letters, **81** (1998), 1322-1325.
- [5] N. A. Gershenfeld and I. L. Chuang, Science **275** (1997), 350-356; D. G. Cory, A. F. Fahmy, and T. F. Havel, Proc. Nat. Acad. Sci. **94** (1997) 1634-1639.
- [6] L. K. Grover, Phys. Rev. Letters, **80** (1998), 4329-4332.
- [7] A. Peres, Phys. Lett. **101A** (1984), 249-250.
- [8] R. Laflamme, C. Miquel, J. P. Paz, and W. Zurek, Phys. Rev. Letters **77** (1996), 198-201; D. P. DiVincenzo and P. W. Shor, Phys. Rev. Letters **77** (1996), 3260-3263; E. M. Raines, R. H. Hardin, P. W. Shor, and N. J. A. Sloane, Phys. Rev. Letters **79** (1997), 953-954; E. Knill, R. Laflamme, and W. Zurek, Science **279** (1998), 342-346; D. Cory, M. Price, W. Maas, E. Knill, R. Laflamme, W. Zurek, T. Havel, and S. Somaroo, Phys. Rev. Letters **81** (1998), 2152-2155.
- [9] H. E. Brandt, Progr. Quantum Electronics, **22** (1998), 257-370.
- [10] P. A. Benioff, Int. Jour. Theoret. Physics **21** (1982), 177-202.
- [11] G. Chen, S. A. Fulling, and M. O. Scully, Los Alamos Archives, quant-ph/9909040.
- [12] M. Boyer, G. Brassard, P. Hoyer, and A. Tapp, Fortschritte Der Physik, **46** (1998), 493-506.
- [13] L. K. Grover, Los Alamos Archives, quant-ph/9912001.
- [14] E. Biham, O. Biham, D. Biron, M. Grassl, and D. Lidar, Phys. Rev. A, **60** (1999), 2742-2745.
- [15] C. Zalka, Phys. Rev. A **60** (1999), 2746-2751.
- [16] E. Farhi and S. Guttmann, Los Alamos Archives, quant-ph/9711035.
- [17] G. L. Long, Y.S. Li, W. L. Zhang, and L. Niu, Physics Letters A, **262** (1999), 27-34.
- [18] J. P. Barnes and W. S. Warren, Phys. Rev. A **60** (1999), 4363-4374.
- [19] S. Lloyd, Phys. Rev. A **61** (2000) 010301(R).
- [20] P. Benioff, Phys. Rev. Letters **78** (1997), 590-593; Phys. Rev. B **55** (1997) 9482-9494.
- [21] L. Viola, S. Lloyd, and E. Knill, Phys. Rev. Letters **83** (1999), 4888-4891; S. Lloyd and J. E. Slotine, Los Alamos Archives quant-ph/9905064.
- [22] W. S. Warren, H. Rabitz, and M Dahleh, Science **259** (1993), 1581- 1589; V. Ramakrishna and H. Rabitz, Phys. Rev. A **51** (1995), 960-966; Phys. Rev. A **54** (1996), 1715-1716; M. Demiralp and H. Rabitz, Phys. Rev. A, **57** (1998), 2420-2425.
- [23] P. Benioff, Phys. Rev. A **59** (1999), 4223-4237.

PHYSICS DIVISION, ARGONNE NATIONAL LABORATORY, ARGONNE, IL 60439

*Current address:* Physics Division, Argonne National Laboratory Argonne, IL 60439

*E-mail address:* pbenioff@anl.gov

PHYSICS DIVISION, ARGONNE NATIONAL LABORATORY, ARGONNE ILLINOIS, 60439

*E-mail address:* pbenioff@anl.gov

