Quantum Formulas: a Lower Bound and Simulation *

Vwani P. Roychowdhury†  Farrokh Vatan‡

Abstract

We show that Nechiporuk’s method [26] for proving lower bounds for Boolean formulas can be extended to the quantum case. This leads to an $\Omega(n^2/\log^2 n)$ lower bound for quantum formulas computing an explicit function. The only known previous explicit lower bound for quantum formulas [27] states that the majority function does not have a linear–size quantum formula. We also show that quantum formulas can be simulated by Boolean circuits of almost the same size.

Key words. quantum formula, lower bound, mixed state, density matrix

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1 Introduction

Computational devices based on quantum physics have attracted much attention lately, and quantum algorithms that perform much faster than their classical counterparts have been developed [12, 21, 22]. To provide a systematic study of the computational power of quantum devices, models similar to those for classical computational devices have been proposed. Deutsch [8] formulated the notion of quantum Turing machine. This approach was further developed by Bernstein and Vazirani [5], and the concept of an efficient universal quantum Turing machine was introduced. As in the case of classical Boolean computation, there is also a quantum model of computation based on circuits (or networks). Yao [27] proved that the quantum circuit model, first introduced by Deutsch [11], is equivalent to the quantum Turing machine model.

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†Electrical Engineering Department, UCLA, Los Angeles, CA 90095 (vwani@ee.ucla.edu).
‡Electrical Engineering Department, UCLA, Los Angeles, CA 90095 (vatan@ee.ucla.edu). Present address: Jet Propulsion Laboratory, California Institute of Technology, 4800 Oak Grove Drive Pasadena, CA 91109 (Farrokh.Vatan@jpl.nasa.gov).
Since every Boolean circuit can be simulated by a quantum circuit, with at most a polynomial factor increase in its size, any nontrivial lower bound for quantum circuits could have far reaching consequences. In classical Boolean circuit theory, all nontrivial lower bounds are for proper subclasses of Boolean circuits such as monotone circuits, formulas, bounded-depth circuits, etc. In the quantum case also it seems that the only hope to prove nontrivial lower bounds is for proper subclasses of quantum circuits. So far the only such known lower bound has been derived by Yao [27] for quantum formulas. The quantum formula is a straightforward generalization of the classical Boolean formula: in both cases, the graph of the circuit is a tree. Yao has proved that the quantum formula size of the majority function \( \text{MAJ}_n \) is not linear, i.e., if \( L(\text{MAJ}_n) \) denotes the minimum quantum formula size of \( \text{MAJ}_n \) then \( \lim_{n \to \infty} \frac{L(\text{MAJ}_n)}{n} = \infty \). This bound is derived from a bound on the quantum communication complexity of Boolean functions.

In this paper, we prove an almost quadratic lower bound for quantum formula size. The key step in the derivation of this lower bound is the extension of Nechiporuk’s method to quantum formulas; for a detailed discussion of Nechiporuk’s method in the Boolean setting see [11, 26]. Nechiporuk’s method has been used in several different areas of Boolean complexity (e.g., see [11] for details). It has also been applied to models where the gates do not take on binary or discrete values, but the input/output map still corresponds to a Boolean function. For example, in [23] this method has been used to get a lower bound for arithmetic and threshold formulas. The challenging part of this method is a step that we shall refer to as “path squeezing” (see §4 for the exact meaning of it). Although in the case of Boolean gates, this part can be solved easily, in the case of analog circuits it is far from obvious (see [23]). For the quantum formulas “path squeezing” becomes even more complicated, because here we should take care of any quantum entanglement and interference phenomena. We show that it is still possible to squeeze a path with arbitrary number of constant inputs to a path with a fixed number of inputs. This leads to a lower bound of \( \Omega(n^2 / \log^2 n) \) on the size of quantum formulas computing a class of explicit functions. For example, we get such a bound for the Element Distinctness function \( \text{ED}_n \). The input of \( \text{ED}_n \), for \( n = 2\ell \log \ell \), is of the form \( (z_1, \ldots, z_\ell) \), where each \( z_j \) is a string of \( 2\log \ell \) bits. Then \( \text{ED}_n(z_1, \ldots, z_\ell) = 1 \) if and only if all these strings are pair wise distinct.

In the end of the paper we compare the powers of quantum formulas and Boolean circuits. Surprisingly, in some sense quantum formulas are not more powerful than Boolean circuits. Any quantum formula of size \( s \) and depth \( d \) can be approximated by a Boolean circuit of size \( O(s \log s \log \log s) \) and depth \( O(d \log \log s) \). Similar results are not known, and most probably are not true, for quantum circuits and other models which are depending on real number parameters (like arithmetic circuits [23]). The key idea for this simulation is that the computation of a quantum formula on an input (which is a pure state in the

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1 There are exponential lower bounds on the time of quantum computation for the black-box model (see, e.g., [6]), but they do not apply to the size of quantum circuits.

2 The value of \( \text{MAJ}_n(x_1, \ldots, x_n) \) is 1 if at least \( \lceil n/2 \rceil \) of inputs are 1.
Hilbert space) can be described as performing a sequence of unitary operations on $4 \times 4$ density matrices of mixed states.

In this paper we use the notation $| \cdot |$ for two different purposes. When $\alpha$ is a complex number, $|\alpha|$ denotes the absolute value of $\alpha$; i.e., $|\alpha| = \sqrt{\alpha \cdot \alpha^*}$. While if $X$ is a set then $|X|$ denotes the cardinality of $X$.

2 Preliminaries

A quantum circuit is defined as a straightforward generalization of acyclic classical (Boolean) circuit (see [10]). For constructing a quantum circuit, we begin with a basis of quantum gates as elementary gates. Each elementary gate $g$ with $d$ inputs represents a unitary operation $U_g \in \mathbf{U}(2^d)$, where $\mathbf{U}(m)$ denotes the group of $m \times m$ unitary complex matrices. The gates are interconnected by quantum “wires”. Each wire represents a quantum bit, qubit, which is a 2-state quantum system represented by a unit vector in $\mathbb{C}^2$. Let $\{|0\rangle, |1\rangle\}$ be the standard orthonormal basis of $\mathbb{C}^2$. The $|0\rangle$ and $|1\rangle$ values of a qubit correspond to the classical Boolean 0 and 1 values, but a qubit can also be in a superposition of the form $\alpha |0\rangle + \beta |1\rangle$, where $\alpha, \beta \in \mathbb{C}$ and $|\alpha|^2 + |\beta|^2 = 1$. Note that the output of such gate, in general, is not a tensor product of its inputs, but an entangled state; e.g., a state like $\frac{1}{\sqrt{2}} |00\rangle + \frac{1}{\sqrt{2}} |11\rangle$ which can not be written as a tensor product.

If the circuit has $m$ inputs, then for each $d$–input gate $g$, the unitary operation $U_g \in \mathbf{U}(2^d)$ can be considered in a natural way as an operator in $\mathbf{U}(2^m)$ by acting as the identity operator on the other $(m - d)$ qubits. Hence, a quantum circuit with $m$ inputs computes a unitary operator in $\mathbf{U}(2^m)$, which is the product of successive unitary operators defined by successive gates.

The size of a quantum circuit $C$, denoted by size($C$), is the number of gates occurring in $C$. The depth of $C$, denoted by depth($C$), is the length of the longest path in $C$ from an input to an output gate.

In this paper, we consider quantum circuits that compute Boolean functions. Consider a quantum circuit $C$ with $m$ inputs. Suppose that $C$ computes the unitary operator $U_C \in \mathbf{U}(2^m)$. We say $C$ computes the Boolean function $f : \{0, 1\}^n \rightarrow \{0, 1\}$ if the following holds. The inputs are labeled by the variables $x_1, x_2, \ldots, x_n$ or the constants $|0\rangle$ or $|1\rangle$ (different inputs may be labeled by the same variable $x_j$). We consider one of the output wires, say the first one, as the output of the circuit. To compute the value of the circuit at $(a_1, \ldots, a_n) \in \{0, 1\}^n$, let the value of each input wire with label $x_j$ be $|a_j\rangle$. These inputs, along with the constant inputs to the circuit, define a unit vector $|\alpha\rangle$ in $\mathbb{C}^{2^n}$. In fact this vector is a standard basis vector of the following form (up to some repetitions and a permutation)

$$|\alpha\rangle = |a_1\rangle \otimes \cdots \otimes |a_n\rangle \otimes |0\rangle \otimes \cdots \otimes |1\rangle.$$ 

The action of the circuit $C$ on the input $|\alpha\rangle$ is the same as $U_C(|\alpha\rangle)$. Note that since $U_C$ is unitary, $\|U_C(|\alpha\rangle)\| = 1$. We decompose the vector $U_C(|\alpha\rangle) \in \mathbb{C}^{2^m}$ with respect to the output qubit. Let the result
Figure 1: Quantum circuits and their computation graphs; the top circuit is not a formula while the bottom one is a formula.

\begin{align*}
U_C(|\alpha\rangle) &= |0\rangle \otimes |A_{0,\alpha}\rangle + |1\rangle \otimes |A_{1,\alpha}\rangle.
\end{align*}

Then we define the probability that \( C \) outputs 1 (on the input \( \alpha \)) as \( p_\alpha = \|A_{1,\alpha}\|^2 \), i.e., the square of the length of \( A_{1,\alpha} \in \mathbb{C}^{2^{m-1}} \). Finally, we say that the quantum circuit \( C \) computes the Boolean function \( f \) if for every \( \alpha \in \{0, 1\}^n \), if \( f(\alpha) = 1 \) then \( p_\alpha > 2/3 \) and if \( f(\alpha) = 0 \) then \( p_\alpha < 1/3 \).

Following Yao [27], we define quantum formulas as a subclass of quantum circuits. A quantum circuit \( C \) is a formula if for every input there is a unique path that connects it to the output qubit. To make this definition more clear we define the computation graph of \( C \), denoted by \( G_C \). The nodes of \( G_C \) correspond to a subset of the gates of \( C \). We start with the output gate of \( C \), i.e., the gate which provides the output qubit, and let it be a node of \( G_C \). Once a node \( v \) belongs to \( G_C \) then all gates in \( C \) that provide inputs to \( v \) are considered as adjacent nodes of \( v \) in \( G_C \). Then \( C \) is a formula if the graph \( G_C \) is a tree. Figure 1 provides examples of quantum circuits of both kinds, i.e., circuits that are also quantum formulas, and circuits that are not formulas.

All circuits that we consider are over some fixed quantum basis. The lower bound does not depend on the basis; the only condition is that the number of inputs (and so the number of outputs) of each gate be bounded by some fixed constant number (this condition is usually considered as part of the definition of a quantum basis). For example, this basis can be the set of all 2–input 2–output quantum gates, and as it is shown in [2], this basis is universal for computation with quantum circuits.
It is well–known that any Boolean circuit can be efficiently simulated by a quantum circuit over a universal basis. Indeed, for this purpose, the 3–bit Toffoli gate is enough (see, e.g., [4, 17]). Similarly, any Boolean formula can be efficiently simulated by a quantum formula using only Toffoli gate or a basis universal for classical computation. In the special case, from [24] it follows that there is a polynomial–size log–depth quantum formula computing the majority function $\text{MAJ}_n$. This fact implies that for quantum formulas over reasonable bases (i.e., universal for classical computation) the threshold probability of correct answer ($\frac{2}{3}$ in the above definition) can be efficiently boosted to a number arbitrarily close to one.

For our proof we also need a Shannon–type result for quantum circuits. Knill [15] has proved several theorems about the quantum circuit complexity of almost all Boolean functions. We will use the following theorem.

**Theorem 2.1** ([15]) *The number of different $n$–variable Boolean functions that can be computed by size $N$ quantum circuits ($n \leq N$) with $d$–input $d$–output elementary gates is at most $2^{cN \log N}$, where $c$ depends only on $d$.***

For the sake of completeness, in Appendix we have provided a proof for a slightly weaker bound. Our approach is different from that in [15] and it seems it is shorter and simpler than the proof in [15]. Although the bound that we get is a little weaker than the bound provided by the above theorem (it is of the form $2^{O(nN)}$), our bound results in the same bound of Theorem 2.1 if $\log(N) = \Omega(n)$ which is true for almost all Boolean functions. Thus our result provides the same bound for the complexity of almost all functions and it is sufficient for the bound we get in this paper.

We also need to consider general orthonormal bases in the space $\mathbb{C}^{2^n}$ other than the standard basis. In the context of quantum physics, we identify the Hilbert space $\mathbb{C}^{2^n}$ as the tensor product space $\bigotimes_{j=1}^{n} \mathbb{C}^2$, and the standard basis consists of the vectors

$$|c_1\rangle \otimes \cdots \otimes |c_n\rangle = |c_1 \cdots c_n\rangle, \ c_j \in \{0, 1\}.$$

**Fact 2.2** *Let $|A_j\rangle \in \mathbb{C}^{2^k}$ and $|B_\ell\rangle \in \mathbb{C}^{2^m}$ be unit vectors (for $j$ and $\ell$ in some index sets). If $|A_j\rangle$ are pair wise orthogonal and $|B_\ell\rangle$ are pair wise orthogonal then the family

$$\left\{ |A_j\rangle \otimes |B_\ell\rangle \in \mathbb{C}^{2^{k+m}} : j, \ell \right\}$$

is an orthonormal set.*

The following lemma, although seemingly obvious, is crucial for the “path squeezing” technique in the proof of the lower bound.
Lemma 2.3  (a) Suppose that $C$ is a subcircuit of a quantum circuit. Let the inputs of $C$ be divided into two disjoint sets of qubits $Q_1$ and $Q_2$. Suppose that each gate of $C$ either acts only on qubits from $Q_1$ or only on qubits from $Q_2$. Then there are subcircuits $C_1$ and $C_2$ such that $C_j$ acts only on qubits from $Q_j$ and the operation of $C$ is the composition of operations of $C_1$ and $C_2$ no matter in which order they act; i.e., $C = C_1 \circ C_2 = C_2 \circ C_1$. So the subcircuit $C$ can be substituted by $C_1$ and $C_2$ (see Figure 2).

(b) Let $C$ be a subcircuit of a quantum circuit with distinct input qubits $q$ and $r_1, \ldots, r_t$. Suppose that only $t$ gates $g_1, \ldots, g_t$ in $C$ act on $q$. Moreover, suppose that each $g_j$ acts on $q$ and $r_j$. Then, w.l.o.g., we can assume that each qubit $r_j$ after entering the gate $g_j$ will not interact with any other qubit until the gate $g_t$ is performed (see Figure 3).

**Proof.** Part (a) is based on the following simple observation. If $M \in \text{U}(2^m)$ and $N \in \text{U}(2^n)$ then

$$M \otimes N = (M \otimes I_n) \circ (I_m \otimes N) = (I_m \otimes N) \circ (M \otimes I_n),$$

where $I_t$ is the identity map in $\text{U}(2^t)$. Note that the inputs of the subcircuit $C$ may be in an entangled state; but to see that the equality $C = C_1 \circ C_2 = C_2 \circ C_1$ holds, it is enough to check this equality for the
standard basis and extend it to the whole space by linearity.

Part (b) follows simply from part (a); as in Figure 4, part (a) can be applied on subcircuit consisting of gates $h_2$ and $h_3$. Note that in this case also input qubits $r_j$ of $g_j$'s may be in an entangled state. Again a linearity argument shows that we have to consider only the case that $r_j$'s are in a product state. 

The above lemma is special case of a more general fact that operations on one part of a bi–partite quantum system do not affect the result of operations on the other part (for more details see, e.g., [18]).

3 A new equivalent definition for quantum formulas

Kitaev [14] has brought to our attention that quantum formulas are equivalent to a model that is very similar to the classical formulas. In this model the inputs and the intermediate results are density matrices. Each gate is a completely positive trace–preserving super–operator, which maps density matrices of a $d$–qubit systems to one–qubit density matrices. The underlying graph, like a classical formula, is a directed tree; i.e., from each input there is a unique path to the output gate. Thus the output of such circuit is a density matrix of a single qubit which provides the probability of the output “0” or “1”. To make the paper self–contained, we first present the definitions of the notions mentioned in this new definition.

By a pure state $|\alpha\rangle$ we mean a unit vector in some Hilbert space $\mathbb{C}^{2^n}$. A mixed state $\{\psi\}$ in $\mathbb{C}^{2^n}$ is a probability distribution on pure states in this Hilbert space. We denote such a mixed state as $\{\psi\} = \{p_k, |\psi_k\rangle\}$, where $p_k \geq 0$ and $\sum_k p_k = 1$. Then $\{\psi\}$ picks the pure state $|\psi_k\rangle$ with probability $p_k$.

The density matrix of a pure state $|\alpha\rangle$ is the matrix $\rho_{|\alpha\rangle}$ of the linear mapping $|\alpha\rangle\langle\alpha|$; i.e, the mapping $|x\rangle \rightarrow \langle\alpha|x\rangle|\alpha\rangle$. So, if $|0\rangle, |1\rangle, \ldots, |2^n-1\rangle$ represent the standard computational basis of $\mathbb{C}^{2^n}$ and $|\alpha\rangle = \sum_k \lambda_k |k\rangle$, then the $(i,j)$ entry of $\rho_{|\alpha\rangle}$ is $\lambda_i \lambda_j^{*}$. The importance of density matrix is that it suffices to characterize the quantum state of the system. Specially, this matrix is enough to find the probabilities of measurements. In general, the result of each measurement can be represented by action of a projection operator $\mathcal{P}$ on the given state $|\alpha\rangle$, where $\mathcal{P}$ is a projection onto some subspace $\mathcal{E}$. Then the probability that the result of the measurement is in the subspace $\mathcal{E}$ is equal to $\text{Tr}(\mathcal{P} \rho_{|\alpha\rangle})$. 

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Figure 4: Changing the order of gates (Lemma 2.3 (b)).
The density matrix of a mixed state $\{\psi\} = \{p_k, |\psi_k\rangle\}$ is defined as

$$\rho_{\{\psi\}} = \sum_k p_k |\psi_k\rangle \langle \psi_k|.$$  

Like the case of pure states, the probability that the result of the measurement is in the subspace $E$ is equal to $\text{Tr}(\mathcal{P} \rho_{\{\psi\}})$.

If the (pure or mixed) state $|\psi\rangle$ can be written as the tensor product $|\phi\rangle \otimes |\chi\rangle$ then the density matrix $\rho_{|\psi\rangle}$ is equal to the tensor (Hadamard) product $\rho_{|\phi\rangle} \otimes \rho_{|\chi\rangle}$.

The next important notion is partial trace. Consider the Hilbert spaces $\mathcal{H}_1 = \mathbb{C}^{2^n}$ and $\mathcal{H}_2 = \mathbb{C}^{2^m}$ and $\mathcal{H} = \mathcal{H}_1 \otimes \mathcal{H}_2$; so $\mathcal{H}$ is isomorphic with $\mathbb{C}^{2^{n+m}}$. Let

$$B_1 = \{ |u_i\rangle : i = 1, \ldots, 2^n \} \quad \text{and} \quad B_2 = \{ |v_j\rangle : j = 1, \ldots, 2^m \}$$

be orthonormal bases for $\mathcal{H}_1$ and $\mathcal{H}_2$, respectively. Then

$$B_1 \otimes B_2 = \{ |u_i\rangle \otimes |v_j\rangle : i = 1, \ldots, 2^n, \ j = 1, \ldots, 2^m \}$$

is a basis for $\mathcal{H}$. Let $\rho$ be the density matrix of a mixed state $|\psi\rangle$ in the space $\mathcal{H}$. It is possible to restrict the state $|\psi\rangle$ to the subspace $\mathcal{H}_1$. The result is a partial trace $\rho|_{\mathcal{H}_1} = \text{Tr}_{\mathcal{H}_2} \rho$ which is density matrix of some mixed state in the subspace $\mathcal{H}_1$. We also say that the subspace $\mathcal{H}_2$ is traced out. The partial trace $\rho|_{\mathcal{H}_1}$ enables us to calculate probabilities of the results of the measurements bearing only on the subspace $\mathcal{H}_1$. We assume that the rows and columns of the matrices $\rho$ and $\rho|_{\mathcal{H}_1}$ are labeled by the vectors in the basis $B_1 \otimes B_2$ and $B_1$, respectively. For example, $\rho(|u_{i_1}\rangle |v_{j_1}\rangle, |u_{i_2}\rangle |v_{j_2}\rangle)$ is the entry of $\rho$ at row labeled by $|u_{i_1}\rangle \otimes |v_{j_1}\rangle$ and the column labeled by $|u_{i_2}\rangle \otimes |v_{j_2}\rangle$. With this notation, the partial trace $\rho|_{\mathcal{H}_1}$ is defined as follows

$$\rho|_{\mathcal{H}_1} (|u_{i_1}\rangle , |u_{i_2}\rangle) = \sum_{j=1}^{2^m} \rho (|u_{i_1}\rangle |v_j\rangle, |u_{i_2}\rangle |v_j\rangle).$$

Again let $E$ be a subspace of $\mathcal{H}_1$. We can identify it with subspace $E \otimes \mathcal{H}_2$ of $\mathcal{H}$. Let $\mathcal{P}: \mathcal{H}_1 \rightarrow E$ be the projection operator associated with $E$. The operator $\mathcal{P}$ can be extended to the whole space $\mathcal{H}$ in a natural way as the operator $\mathcal{P} \otimes \text{Id}_{\mathcal{H}_2}$, where $\text{Id}_{\mathcal{H}_2}$ is the identity operator on $\mathcal{H}_2$. Then the probability that the result of the measurement is in the subspace $E$ is equal to $\text{Tr}(\mathcal{P} \rho|_{\mathcal{H}_1})$. (For more details on density matrices of mixed states and the partial trace see, e.g., [8].)

Let $C$ be a quantum circuit. For inputs of $C$ it is possible to consider mixed states along with pure states. Toward this end, each input is substituted by its density matrix, and each gate $g$ of $C$ by a superoperator $\bar{g}$ that maps density matrices to density matrices. In fact, if the unitary operator of the gate $g$ is $U$, then the action of $\bar{g}$ on the density matrix $\rho$ is as follows:

$$\bar{g}(\rho) = g \circ \rho = U \rho U^\dagger. \quad (1)$$
Lemma 3.1 ([1]) If the gates $g_1$ and $g_2$ operate on disjoint sets of qubits, then for any density matrix $\rho$ we have $g_1 \circ g_2 \circ \rho = g_2 \circ g_1 \circ \rho$.

First we show that every quantum formula is equivalent to a circuit based on this new definition. Let $F$ be a quantum formula on a basis of $d$–bit gates. Construct a circuit $C$ from $F$ by the following transformations. In each gate $g$, performing the unitary operation $U \in U(2^d)$, keep the only output which is connected to the output and substitute the operator $g$ by the super–operator $[g] = \text{Tr}_{\tilde{H}} \circ \tilde{g}$, where $\tilde{H}$ is the $(d-1)$–dimensional subspace spanned by the qubits removed from the output of this gate. The fact that the circuit $C$ computes the same function as the formula $F$ follows from Lemma 2.3. Thus the underlying graph of the circuit $C$ is the same as the computation tree of the formula $F$, where the node corresponded with the gate $g$ computes the super–operator $[g]$.

Let now $C$ be a circuit based on this new definition. We construct a quantum formula $F$ from $C$ by simply substituting each gate of $C$, computing the super–operator $T$, by a $(d+2)$–input $(d+2)$–output unitary gate $U$, only one output of this gate is connected to the next gate and the other outputs never interact with any other qubit. So $F$ satisfies our original definition of quantum formula. The only thing remains is to show how we can choose the unitary operators $U$ such that the formula $F$ computes the same Boolean function as $C$. The following theorem guarantees the existence of the correct operator $U$, for each gate of $C$. Here $L(\mathcal{H})$ is the space of linear operators on the Hilbert space $\mathcal{H}$ and for unitary operator $U$ on $\mathcal{H}$, the operator $O_U \in L(\mathcal{H})$ is defined as $O_U(M) = UMU^\dagger$.

**Theorem 3.2 ([1, 13, 16, 20])** Suppose that $T : L(\mathcal{H}_1) \longrightarrow L(\mathcal{H}_2)$ is a trace–preserving and completely positive super–operator. Then there are Hilbert spaces $\mathcal{G}_1$ and $\mathcal{G}_2$, where $\dim (\mathcal{G}_1) = (\dim (\mathcal{H}_2))^2$ and $\dim (\mathcal{G}_2) = \dim (\mathcal{H}_1) \cdot \dim (\mathcal{H}_2)$, and there is a unitary operator $U : \mathcal{H}_1 \otimes \mathcal{G}_1 \longrightarrow \mathcal{H}_2 \otimes \mathcal{G}_2$ such that $T = \text{Tr}_{\mathcal{G}_2} \circ O_U$.

We would like to mention that from now on it might be more useful to accept the new modified definition as the standard one for quantum formulas in the literature.

4 The lower bound

Let $f(x_1, \ldots, x_n)$ be a Boolean function. Let $X = \{x_1, \ldots, x_n\}$ be the set of the input variables. Consider a partition $\{S_1, \ldots, S_k\}$ of $X$; i.e.,

$$X = \bigcup_{j=1}^{k} S_j \quad \text{and} \quad S_{j_1} \cap S_{j_2} = \emptyset, \quad \text{for} \quad j_1 \neq j_2.$$ 

Let $n_j = |S_j|$, for $j = 1, \ldots, k$. Let $\mathcal{F}_j$ be the set of all subfunctions of $f$ on $S_j$ obtained by fixing the variables outside $S_j$ in all possible ways. We denote the cardinality of $\mathcal{F}_j$ by $\sigma_j$. 

}\begin{align*}
L(\mathcal{H}_1) \longrightarrow L(\mathcal{H}_2) & \text{ is a trace–preserving and completely positive super–operator. Then there are Hilbert spaces } \mathcal{G}_1 \text{ and } \mathcal{G}_2, \text{ where } \\
\dim (\mathcal{G}_1) & = (\dim (\mathcal{H}_2))^2 \text{ and } \\
\dim (\mathcal{G}_2) & = \dim (\mathcal{H}_1) \cdot \dim (\mathcal{H}_2), \text{ and there is a unitary operator } U : \mathcal{H}_1 \otimes \mathcal{G}_1 \longrightarrow \mathcal{H}_2 \otimes \mathcal{G}_2 \text{ such that } \\
T & = \text{Tr}_{\mathcal{G}_2} \circ O_U. 
\end{align*}
As an example, we compute the above parameters for the Element Distinctness function ED\(_n\) (see [6]). Let \(n = 2\ell \log \ell\) (so \(\ell = \Omega(n/\log n)\)) and divide the \(n\) inputs of the function into \(\ell\) strings each of \(2\log \ell\) bits. Then the value of ED\(_n\) is 1 if and only if these \(\ell\) strings are pair wise distinct. We consider the partition \((S_1, \ldots, S_\ell)\) such that each \(S_j\) contains all variables of the same string. Thus \(n_j = |S_j| = 2\log \ell\). Each string in \(S_j\) represents an integer from the set \(\{0, 1, \ldots, \ell^2 - 1\}\). The function ED\(_n\) is symmetric with respect to \(S_j\)'s; so \(|F_j| = |F_j'|\). To estimate \(|F_1|\), note that if the strings \((z_2, \ldots, z_\ell)\) in \(S_2, \ldots, S_\ell\) represent distinct integers then the corresponding subfunction is different from any subfunction corresponding to any other string. So \(\sigma_j = |F_1| \geq (\ell^2/\ell - 1) > \ell^{\ell-1}\).

**Theorem 4.1** Every quantum formula computing \(f\) has size

\[
\Omega \left( \sum_{1 \leq j \leq k} \frac{\log(\sigma_j)}{\log \log(\sigma_j)} \right).
\]

**Proof.** We give a proof for any basis consisting of 2–input 2–output quantum gates. The proof for bases with more than two inputs is a simple generalization of this proof.

Let \(F\) be a formula computing \(f\). Let \(\Sigma_j\) be the set of input wires of \(F\) labeled by a variable from \(S_j\), and let \(s_j = |\Sigma_j|\). Then

\[
\text{size}(F) = \Omega \left( \sum_{1 \leq j \leq k} s_j \right).
\]

We want to consider the formulas obtained from \(F\) by letting the input variables not in \(\Sigma_j\) to some constant value \(|0\rangle\) or \(|1\rangle\). In this regard, let \(P_j\) be the set of all paths from an input wire in \(\Sigma_j\) to the output of \(F\). Finally, let \(G_j\) be the set of gates of \(F\) where two paths from \(P_j\) intersect. Then \(|G_j| \leq s_j\).

Let \(\tau\) be an assignment of \(|0\rangle\) or \(|1\rangle\) to the input variable wires *not* in \(\Sigma_j\). We denote the resulting formula by \(F_\tau\). Thus \(F_\tau\) computes a Boolean function \(f_\tau : \{0, 1\}^{n_j} \rightarrow \{0, 1\}\) which is a subfunction of \(f\) and a member of \(F_j\). Consider a path

\[
\pi = (g_1, g_2, \ldots, g_m), \quad m > 2,
\]

in \(F_\tau\), where \(g_1\) is an input wire or a gate in \(G_j\), \(g_m\) is a gate in \(G_j\) or the output wire of \(F\), and \(g_\ell \notin G_j\) for \(1 < \ell < m\).

To show how we can squeeze paths like \(\pi\) (this is the essence of the Nechiporuk’s method), we introduce the following notations. We consider a natural ordering \(\gamma_1, \gamma_2, \ldots, \gamma_t\) on the gates of the formula \(F_\tau\), and regard \(F_\tau\) as a computation in \(t\) steps where at step \(\ell\) the corresponding gate \(\gamma_\ell\) is performed. We say two qubits \(q_1\) and \(q_2\) are *strong companions* of each other at step \(\ell\) if there is a gate \(\gamma_j\) such that \(j \leq \ell\)
Figure 5: The qubits $q_1$ and $q_2$ are strong companions at step $\ell$, the qubits $q_2$ and $q_3$ are companions at step $\ell + 2$.

and $q_1$ and $q_2$ are inputs of $\gamma_j$. We say qubits $q_1$ and $q_2$ are companions of each other at step $\ell$ if there exists a sequence $r_1, r_2, \ldots, r_p$ of qubits such that $r_1 = q_1$, $r_p = q_2$, and $r_j$ and $r_{j+1}$ (for $1 \leq j \leq p - 1$) are strong companions of each other at step $\ell$ (see Figure 5). If $q_1$ and $q_2$ are companions at step $\ell$ then they are also companions at any step after $\ell$. For a gate $g = \gamma_k$, we define the set of companions of $g$ as the union of all companions of input qubits of $g$ at step $k$.

Suppose that in the path (3) $g_1 = \gamma_j_0$, $g_m = \gamma_j_1$, the inputs of $g_1$ are $q_0$ and $q_1$, the output of $\gamma_j_0$ from the path (3) is the qubit $q_0$, and the input of $\gamma_j_1$ not from the path (3) is the qubit $q_2$ (see Figure 5). Note that $q_0$ is the companion of $q_2$ at step $j_1$. Let $Q_\pi$ be the union of all sets of companions of $g_2, \ldots, g_{m-1}$ minus $q_0$ and $q_1$ and their companions at step $j_1$. Let $C_0$ be the circuit defined by the gates $g_1, \ldots, g_{m-1}$ from the path (3). Suppose that $|Q_\pi| = v$ and consider $C_0$ as an operation acting on $\mathcal{H} = \mathbb{C}^2 \otimes \mathbb{C}^2 \otimes \mathbb{C}^2^v$. To study the action of the subcircuit $C_0$, it is enough to consider the action of $C_0$ on the computational basis vectors of the space $\mathcal{H}$. Therefore, while the inputs of $C_0$ as a subcircuit of $F_\tau$ are in general entangled states, we only have to study the action of $C_0$ on the computational basis vectors which are product states. We label the inputs $|\alpha_0\rangle \otimes |\alpha_1\rangle \otimes |\alpha\rangle \in \mathcal{H}$ of $C_0$ in such a way that when $C_0$ acts as a subformula of $F_\tau$ then $|\alpha_0\rangle$,
\[ |\alpha_0\rangle \otimes |\alpha_1\rangle \otimes |0\ldots0\rangle \longrightarrow \sum_{c_0,c_1 \in \{0,1\}} |c_0\rangle \otimes |c_1\rangle \otimes |A^0_{c_0,c_1}\rangle, \quad (4) \]

where \( \alpha_0, \alpha_1 \in \{0,1\} \), and \( |A^0_{c_0,c_1}\rangle \in \mathbb{C}^{2^n} \) may be not a unit vector. Let \( A_\pi \subseteq \mathbb{C}^{2^n} \) be the vector space spanned by \( |A^0_{c_0,c_1}\rangle \), for \( \alpha_0, \alpha_1, c_0, c_1 \in \{0,1\} \) and \( d = \dim(A_\pi) \). Then \( 1 \leq d \leq 16 \). Let \( |A^7_1\rangle, \ldots, |A^7_d\rangle \) be an orthonormal basis for \( A_\pi \). Then we can rewrite (4) as follows

\[ |\alpha_0\rangle \otimes |\alpha_1\rangle \otimes |0\ldots0\rangle \longrightarrow \sum_{c_0,c_1 \in \{0,1\}} \sum_{1 \leq j \leq d} \lambda^0_{j,c_0,c_1} |c_0\rangle \otimes |c_1\rangle \otimes |A^j_j\rangle. \quad (5) \]

Let \( \mathcal{M}_\pi \) be the set of those unitary operations that are performed after one of the gates \( g_1, \ldots, g_{m-1} \) on some qubits in \( Q_\pi \) before the step \( j_1 \). Since qubits in \( Q_\pi \) do not interact with any other path of the form (3), by Lemma 2.3 (b), we can postpone all operations in \( \mathcal{M}_\pi \) after we computed the output of \( g_m \). Let \( \pi_1, \ldots, \pi_k \) be a natural ordering on the paths like (3) on all paths in \( P_j \) (i.e., the last gate of \( \pi_j+1 \) is not performed before the last gate of \( \pi_j \)). Consider the sets of postponed operations \( \mathcal{M}_{\pi_1}, \ldots, \mathcal{M}_{\pi_k} \). Once again Lemma 2.3 implies that we can postpone operations in \( \mathcal{M}_{\pi_1} \) after the last gate of \( \pi_2 \); then we can postpone operations in \( \mathcal{M}_{\pi_1} \) and \( \mathcal{M}_{\pi_2} \) after the last gate of \( \pi_3 \), and so on. Repeating this argument shows that we can postpone all operations in \( \mathcal{M}_{\pi_1}, \ldots, \mathcal{M}_{\pi_k} \) after we compute the output qubit. In this way, the state of the output qubit, before the postponed operations \( \mathcal{M}_{\pi_1}, \ldots, \mathcal{M}_{\pi_k} \) are applied, is of the form

\[ |0\rangle \otimes |M\rangle + |1\rangle \otimes |N\rangle, \quad (6) \]

where the first qubit is the output qubit and \( |M\rangle \) and \( |N\rangle \) are superpositions of tensor products of orthonormal vectors \( |A^7_j\rangle \) used in (5). By Fact 2.2 these tensor products of the vectors \( |A^7_j\rangle \) are unit vectors and pair wise orthogonal. The unitary operations in the sets \( \mathcal{M}_{\pi_j} \) (for paths \( \pi_j \) of the form (3)), which are postponed to the end, do not change the lengths of \( |M\rangle \) and \( |N\rangle \). Thus, as far as the computation of the Boolean function \( f_\pi \) is concerned, we can ignore all the postponed unitary operations. For this reason we construct the circuit \( \overline{F_\pi} \) from the formula \( F_\pi \) by eliminating all postponed operations in \( \mathcal{M}_{\pi_j} \), substituting for each path \( \pi_j \) of the form (3) the companion qubits in \( Q_{\pi_j} \) by four new qubits, and the unitary operation (5) by the operation defined as

\[ |\alpha_0\rangle \otimes |\alpha_1\rangle \otimes |0000\rangle \longrightarrow \sum_{c_0,c_1 \in \{0,1\}} \sum_{0 \leq j \leq 15} \lambda^0_{j,c_0,c_1} |c_0\rangle \otimes |c_1\rangle \otimes |j\rangle. \quad (7) \]
The output of the circuit $F_\tau$, instead of (6), is of the form
\[ |0\rangle \otimes |M'\rangle + |1\rangle \otimes |N'\rangle, \tag{8} \]
where $\|M\| = \|M'\|$ and $\|N\| = \|N'\|$. So the circuit $F_\tau$ computes $f_\tau$. Moreover,
\[ \text{size}(F_\tau) = O(s_j), \]
and for another assignment $\tau'$, the corresponding circuit $F_{\tau'}$ differs from $F_\tau$ only at unitary operations defined by (6).

The above discussion implies that $\sigma_j$, the number of subfunctions on $S_j$, is at most the number of different Boolean functions computed by size $O(s_j)$ quantum circuits. Therefore, by Theorem 2.1, we get
\[ \sigma_j \leq 2^{O(s_j \log s_j)}. \]
So $s_j = \Omega(\log(\sigma_j)/\log \log(\sigma_j))$. Now the theorem follows from (2).

We would like to mention that the fact that a path like (3) can be squeezed to a path of constant length is a special case of the general property of super–operators stated in Theorem 3.2.

To apply the general bound of the above theorem, we could consider any of the several explicit functions used in the case of Boolean formulas (see [11, 26]). As we mentioned in the beginning of this section, we consider the Element Distinctness function $\text{ED}_n$. For this function $\sigma_j > \ell^{\ell-1}$, where $\ell = \Omega(n/ \log n)$. Therefore, we get the lower bound $\Omega(\ell^2) = \Omega(n^2/ \log^2 n)$ for the formula size.

**Theorem 4.2** Any quantum formula computing $\text{ED}_n$ has size $\Omega(n^2/ \log^2 n)$.

## 5 Quantum formulas vs. Boolean circuits

In this section we show that quantum formulas are not more powerful than Boolean circuits. So as a model of computation, their strength lies between Boolean formulas and Boolean circuits.

Following the idea developed in Section 3, we consider a quantum formula as a quantum circuit operating on mixed states. For the details of quantum circuits with mixed states see [1, 13]. Before we start the proof of the main result of this section, we need to see how we can bound errors in quantum circuits with mixed states. Toward this end we need a suitable norm on super–operators. Each super–operator $T$ which maps density matrices to density matrices is a linear mapping of the form $L(\mathcal{H}_1) \longrightarrow L(\mathcal{H}_2)$, where $\mathcal{H}_1$ and $\mathcal{H}_2$ are finite–dimensional Hilbert spaces and $L(\mathcal{H}_j)$ is the set of linear operators on $\mathcal{H}_j$. Note that $L(\mathcal{H}_j)$ itself is a linear space. Let $\mathcal{H}$ be an $m$–dimensional Hilbert space. There are several norms on the
space $\mathbf{L}(\mathcal{H})$, of which we need the following ones. Let $A \in \mathbf{L}(\mathcal{H})$. We identify $A$ with its $m \times m$ matrix $(a_{ij})$. The first norm is

$$M(A) = m \max_{i,j} |a_{ij}|.$$ 

The usual norm is defined as

$$\|A\| = \sup_{|x| \neq 0} \frac{\|A|x\|}{\|x\|} = \max \left\{ \sqrt{\lambda} : \lambda \in \text{Spec}(A^\dagger A) \right\},$$

where $\text{Spec}(M)$ is the spectrum of the matrix $M$; i.e., the set of the eigenvalues of $M$. The other norm is the trace norm:

$$\|A\|_{\text{Tr}} = \sum_{\lambda \in \text{Spec}(A^\dagger A)} \sqrt{\lambda}.$$ 

We need the next norm $\|\cdot\|_*$ to define another norm: let $T$ be a linear operator that maps matrices to matrices; i.e., $T \in \mathbf{L}(\mathbf{L}(\mathcal{H}))$, then

$$\|T\|_* = \sup_{A \neq 0} \frac{\|TA\|_{\text{Tr}}}{\|A\|_{\text{Tr}}}. $$

The last norm we consider is the diamond norm, defined in [13] and also in [1]. To define this norm, we consider a Hilbert space $\mathcal{G}$ such that $\dim(\mathcal{G}) \geq \dim(\mathcal{H})$ and we let

$$\|T\|_\diamond = \|T \otimes I_\mathcal{G}\|_*,$$

where $I_\mathcal{G}$ is the identity operator on $\mathcal{G}$. The followings are the basic properties of these norms.

(i) $\frac{1}{m} M(A) \leq \|A\| \leq M(A)$.

(ii) $\|A\|_{\text{Tr}} \leq m \|A\|.$

(iii) $\|T(\rho)\|_{\text{Tr}} \leq \|T\|_\diamond \|\rho\|_{\text{Tr}}$, for the density matrix $\rho$.

(iv) $\|TR\|_\diamond \leq \|T\|_\diamond \|R\|_\diamond$.

(v) $\|T \otimes R\|_\diamond = \|T\|_\diamond \|R\|_\diamond$.

(vi) If $T = \tilde{g}$, for some quantum gate $g$, or $T = \text{Tr}_\mathcal{F}$, then $\|T\|_\diamond = 1$.

The properties (iii)–(vi) are proved in [1, 13].

For any operator $V \in \mathbf{L}(\mathcal{H})$ we define the operator $\mathcal{O}_V \in \mathbf{L}(\mathbf{L}(\mathcal{H}))$ as

$$\mathcal{O}_V(M) = V M V^\dagger, \quad M \in \mathbf{L}(\mathcal{H}).$$

(9)

In [1, 13] it is proved that $\|\mathcal{O}_V - \mathcal{O}_W\|_\diamond \leq 2 \|V - W\|$ if $\|V\| \leq 1$ and $\|W\| \leq 1$. We need the following general form of this inequality.
Lemma 5.1 Let \( \dim(\mathcal{H}) = m \). For any \( V, W \in L(\mathcal{H}) \) we have
\[
\|\mathcal{O}_V - \mathcal{O}_W\|_\infty \leq 2m \|V - W\| \min(\|V\|, \|W\|) + m \|V - W\|^2.
\]

Proof. We have (for \( A \in L(\mathcal{H} \otimes \mathcal{H}) \))
\[
\|\mathcal{O}_V - \mathcal{O}_W\|_\infty = \sup_{A \neq 0} \left\| (\mathcal{O}_V \otimes I_{3\mathcal{H}}) A - (\mathcal{O}_W \otimes I_{3\mathcal{H}}) A \right\|_{\text{Tr}} / \|A\|_{\text{Tr}}
\]
\[
= \sup_{A \neq 0} \left\| (V \otimes I_{3\mathcal{H}}) A (V^\dagger \otimes I_{3\mathcal{H}}) - (W \otimes I_{3\mathcal{H}}) A (W^\dagger \otimes I_{3\mathcal{H}}) \right\|_{\text{Tr}} / \|A\|_{\text{Tr}}
\]
\[
= \sup_{A \neq 0} \left\| (V \otimes I_{3\mathcal{H}}) A (V^\dagger \otimes I_{3\mathcal{H}}) - ((V + (W - V)) \otimes I_{3\mathcal{H}}) A ((V^\dagger + (W^\dagger - V)) \otimes I_{3\mathcal{H}}) \right\|_{\text{Tr}} / \|A\|_{\text{Tr}}
\]
\[
\leq \sup_{A \neq 0} \left\| (V \otimes I_{3\mathcal{H}}) A ((W^\dagger - V^\dagger) \otimes I_{3\mathcal{H}}) \right\|_{\text{Tr}} / \|A\|_{\text{Tr}} + \sup_{A \neq 0} \left\| ((W - V) \otimes I_{3\mathcal{H}}) A (V^\dagger \otimes I_{3\mathcal{H}}) \right\|_{\text{Tr}} / \|A\|_{\text{Tr}}.
\]

Since \( MN \leq \|M\| \cdot \|N\|, \|M\| \leq \|M\|_{\text{Tr}} \leq m \|M\|, \|M \otimes N\| = \|M\| \cdot \|N\|, \|M^\dagger\| = \|M\|, \text{ and} \|I_{3\mathcal{H}}\| = 1 \), it follows that
\[
\left\| (V \otimes I_{3\mathcal{H}}) A ((W^\dagger - V^\dagger) \otimes I_{3\mathcal{H}}) \right\|_{\text{Tr}} \leq m \left\| (V \otimes I_{3\mathcal{H}}) A ((W^\dagger - V^\dagger) \otimes I_{3\mathcal{H}}) \right\|
\]
\[
\leq m \|V \otimes I_{3\mathcal{H}}\| \cdot \|A\| \cdot \left\| (W^\dagger - V^\dagger) \otimes I_{3\mathcal{H}} \right\|
\]
\[
\leq m \|V\| \cdot \|A\|_{\text{Tr}} \cdot \|W - V\|.
\]

By applying a similar reduction to the other terms of (10), we drive the following inequality
\[
\|\mathcal{O}_V - \mathcal{O}_W\|_\infty \leq 2m \|V - W\| \cdot \|V\| + m \|V - W\|^2.
\]

We can also drive a similar inequality with \(\|V\|\) substituted by \(\|W\|\). This completes the proof.

We say two \(n \times m\) matrices \(A = (a_{ij})\) and \(B = (b_{ij})\) are \(\delta\)-close to each other if \(|a_{ij} - b_{ij}| \leq \delta\), for every \(1 \leq i \leq n\) and \(1 \leq j \leq m\). If the \(n \times m\) matrices \(A\) and \(B\) are \(\delta\)-close to each other then
\[
\|A - B\| \leq M(A - B) \leq m \delta.
\]

The following theorem formalizes the general form of the error bound for quantum circuits when approximating the unitary operator of each gate. This theorem is actually a generalization of a weaker
theorem which has appeared in several papers (see, e.g., [1, 5, 13]). We need this generalization because once we substitute any unitary gate $S$ of the original quantum circuit by some approximated gate $T$, in general we do not know whether $\|T\| \leq 1$ or not (this is the assumption of the weaker version of this theorem).

**Theorem 5.2** Let $S_j, T_j \in L \left( C^{2^d} \right)$, $1 \leq j \leq \ell$, be defined as $S_j = \mathcal{O}_{U_j}$ and $T_j = \mathcal{O}_{V_j}$, where $U_j \in U(2^d)$ is unitary and $V_j$ is $\delta$–close to $U_j$. Then

$$\|S_\ell \cdots S_3 S_2 S_1 - T_\ell \cdots T_3 T_2 T_1\|_o \leq e^{\eta(d,\delta)\ell} - 1,$$

where $\eta(d, \delta) = 2^{2d+1}\delta (1 + 2^d \delta)$.

**Proof.** First note that

$$\|S_j - T_j\|_o = \|\mathcal{O}_{U_j} - \mathcal{O}_{V_j}\|_o$$

$$\leq 2^{d+1} \|U_j - V_j\| (1 + \|U_j - V_j\|) \quad \text{by Lemma 5.1}$$

$$\leq 2^{2d+1}\delta (1 + 2^d \delta) \quad \text{by (11)}$$

$$= \eta(d, \delta);$$

and, by (vi),

$$\|T_j\|_o \leq \|S_j\|_o + \|S_j - T_j\|_o \leq 1 + \eta(d, \delta). \quad (12)$$

Also we have the following simple inequality

$$\|M_2 M_1 - N_2 N_1\|_o = \|M_2(M_1 - N_1) - (M_2 - N_2)N_1\|_o$$

$$\leq \|M_2\|_o \|M_1 - N_1\|_o + \|N_1\|_o \|M_2 - N_2\|_o \quad (13)$$

Now, by repeated applications of (12) and (13), we have

$$\|S_\ell \cdots S_3 S_2 S_1 - T_\ell \cdots T_3 T_2 T_1\|_o \leq \|S_\ell \cdots S_3 S_2\|_o \|S_1 - T_1\|_o +$$

$$\|T_1\|_o \|S_\ell \cdots S_3 S_2 - T_\ell \cdots T_3 T_2\|_o$$

$$\leq \eta(d, \delta) + (1 + \eta(d, \delta)) \|S_\ell \cdots S_3 S_2 - T_\ell \cdots T_3 T_2\|_o$$

$$\leq \eta(d, \delta) + \eta(d, \delta)(1 + \eta(d, \delta)) +$$

$$(1 + \eta(d, \delta))^2 \|S_\ell \cdots S_3 - T_\ell \cdots T_3\|_o$$

$$\vdots$$

$$\leq \eta(d, \delta) \sum_{j=0}^{\ell-1} (1 + \eta(d, \delta))^j$$

$$= (1 + \eta(d, \delta))^\ell - 1$$

$$\leq e^{\eta(d,\delta)\ell} - 1.$$
The following theorem is the immediate consequence of the above theorem. Note that for a gate $g$ the super–operator $\tilde{g}$ is defined by \([1]\).

**Theorem 5.3** Let $C$ be a quantum circuit composed of the gates $g_1, \ldots, g_s$. Suppose that each $g_j$ is a $d$–bit gate computing the unitary operator $U_j \in U(2^d)$. For each $1 \leq j \leq s$, let $V_j \in L\left(\mathbb{C}^{2^d}\right)$ be a $\delta$–close matrix to $U_j$. Let $T_j \in L\left(L\left(\mathbb{C}^{2^d}\right)\right)$ be defined as $T_j = O_{V_j}$. For any input density matrix $\rho_0$, let

$$
\psi = (\tilde{g}_s \otimes I_{\mathcal{H}_s}) \circ \cdots \circ (\tilde{g}_2 \otimes I_{\mathcal{H}_2}) \circ (\tilde{g}_1 \otimes I_{\mathcal{H}_1}) \rho_0
$$

be the output of $C$, where $\mathcal{H}_j$ is the Hilbert space generated by the qubits not involved with the gate $g_j$. Also, let

$$
\zeta = (T_s \otimes I_{\mathcal{H}_s}) \cdots (T_2 \otimes I_{\mathcal{H}_2})(T_1 \otimes I_{\mathcal{H}_1}) \rho_0
$$

be the approximated output of C. Then

$$
\|\psi - \zeta\|_{\text{Tr}} \leq \left(e^{\eta(d,\delta)s} - 1\right) \|\rho_0\|_{\text{Tr}},
$$

where $\eta(d,\delta) = 2^{2d+1}\delta(1 + 2^d\delta)$.

**Theorem 5.4** Let $\mathcal{B}$ be a quantum basis. Then each quantum formula of size $\ell$ and depth $d$ over the basis $\mathcal{B}$ can be simulated with error at most $\varepsilon$ by a Boolean circuit of size $O(\ell \mu \log \mu \log \log \mu)$ and depth $O(d \log \mu)$, where $\mu = \lceil \log \ell - \log \varepsilon \rceil$.

**Proof.** The basic idea of the simulation is to look at the behavior of a quantum formula as a quantum circuit acting on density matrices of mixed states. We assume, w.l.o.g., that each gate in the basis $\mathcal{B}$ is a 2–bit gate.

Consider a quantum formula $F$ over the basis $\mathcal{B}$; suppose that $F$ has $t$ inputs (constant or variable) and computes the Boolean function $f : \{0, 1\}^n \rightarrow \{0, 1\}$. We show that there is a Boolean circuit $\mathcal{C}$ that for any input $\alpha = (a_1, \ldots, a_n) \in \{0, 1\}^n$ simulates the action of $F$ on $\alpha$. Let $|\alpha\rangle = |0\rangle \otimes |A_0\rangle + |1\rangle \otimes |A_1\rangle$ be the output of $F$ on the input $\alpha$. Suppose that the first qubit is the output bit. If we trace out the non–output bits of $|\alpha\rangle$, the result is a $2 \times 2$ density matrix $\rho_{\text{final}} = \rho|_{\alpha}$. From $\rho_{\text{final}}$ it is easy to calculate the probability of acceptance of $F$. The formula structure of $F$ allows us to calculate the density matrix $\rho_{\text{final}}$ without going to the $2^t$ dimensional space. The Boolean circuit $\mathcal{C}$ finds the density matrix $\rho_{\text{final}}$ by simulating the gates of $F$ step by step.

Since the trace norm of a density matrix is equal to its trace, it follows that $\|\rho_0\|_{\text{Tr}} = 1$, where $\rho_0$ is the density matrix of the input.
Now the gates of $\mathcal{F}$ are no longer acting on pure states, but they are acting on mixed states. If the input of a gate $g_j$, (performing the unitary operation $U_j$) is the $4 \times 4$ density matrix $\rho$ then the output is the density matrix $\rho' = U_j \rho U_j^\dagger$. Of the two output bits $q_1$ and $q_2$ of this gate only one, say $q_1$, is connected to the output bit of $\mathcal{F}$. So we trace out the system representing $q_2$ and consider the new density matrix $\rho_{q_1} = \Tr_{q_2} \rho'$ for $q_1$. By repeating this process for each gate of $\mathcal{F}$ we finally get the desired density matrix $\rho_{\text{final}}$. The correctness of this process follows from Lemma 3.1.

The Boolean circuit $\mathcal{C}$ can simulate the calculations of these density matrices $\rho_{q_1}$. The only problem for this simulation is the proper approximation of the entries of unitary matrices $U_j$. If we substitute each entry of $U_j$ by its first $\mu = -\lceil \log_2 \delta \rceil$ bits, then we get a matrix that is $\delta$–close to $U_j$. Let $\tilde{\mathcal{F}}$ be the resulting formula and $\tilde{\rho}_{\text{final}}$ be the output of $\tilde{\mathcal{F}}$. Then, by Theorem 5.3, $\| \rho_{\text{final}} - \tilde{\rho}_{\text{final}} \|_{\text{Tr}} \leq e^{\eta(d,\delta)\ell} - 1$. So if $\delta = O\left(\frac{\varepsilon}{\ell}\right)$, i.e., $\mu = O(\log \ell - \log \varepsilon)$, then the simulation of $\mathcal{F}$ by $\tilde{\mathcal{F}}$ has at most $\varepsilon$ error. The theorem now follows from this fact that addition and multiplication of $m$ bits numbers can be carried out by Boolean circuits of size $O(m \log m \log \log m)$ and depth $O(\log m)$ (see [19, 26]).

Why does this proof not provide a Boolean formula instead of Boolean circuit? The reason is that to calculate $\rho' = U_j \rho U_j^\dagger$, we need 4 copies of each entry of $\rho$. Thus the fan–out of the gates in the Boolean circuit obtained from the formula $\tilde{\mathcal{F}}$ is 4. This means that the Boolean formula equivalent to this Boolean circuit, in general, has size exponential in $\ell$; this size is at least $\Omega(\ell^3)$, if the graph of $\mathcal{F}$ is a full binary tree.

6 Concluding Remarks

We have extended a classical technique for proving lower bound for Boolean formula size to quantum formulas. The difficult part was to effectively deal with the phenomenon of entanglement of qubits. While we have been successful in extending a classical technique to the quantum case, the challenges encountered indicate that in general the problem of extending methods of Boolean case to the quantum case may not have simple solutions. For example, even the seemingly simple issue of the exact relationship between quantum formulas and quantum circuits has not been resolved. In the Boolean case, simulation of circuits by formulas is a simple fact, but in the quantum case it is not clear whether every quantum circuit can be simulated by a quantum formula. In particular, it is not clear that in the process of going from quantum circuits to formulas, how we can modify the underlying entanglement of qubits while keeping the probability of reaching to the final answer the same. We were also able to show that it is possible to simulate quantum formulas with Boolean circuits of almost the same size. It does not seem that Boolean formulas could efficiently simulate their quantum counterparts. So evidently quantum formulas, as a model of computation, are more powerful than Boolean formulas and less powerful than Boolean circuits. A better
understanding of the relations between these models remains a challenging problem.

7 Appendix: Counting the number of Boolean functions computed by quantum circuits of a given size

In this appendix we prove the following upper bound.

**Theorem 7.1** The number of different $n$–variable Boolean functions that can be computed by size $N$ quantum circuits ($n \leq N$) with $d$–input $d$–output elementary gates (for some constant $d$) is at most $2^{O(nN) + O(N \log N)}$.

Our proof is based on Warren’s bound on the number of different sign–assignments to real polynomials [25]. We begin with some necessary notations.

Let $P_1(x_1, \ldots, x_t), \ldots, P_m(x_1, \ldots, x_t)$ be real polynomials. A sign–assignment to these polynomials is a system of inequalities

$$P_1(x_1, \ldots, x_t) \Delta_1 0, \ldots, P_m(x_1, \ldots, x_t) \Delta_m 0,$$

where each $\Delta_j$ is either “$<$” or “$>$”. The sign–assignment (14) is called consistent if this system has a solution in $\mathbb{R}^t$.

**Theorem 7.2** (Warren [25]) Let $P_1(x_1, \ldots, x_t), \ldots, P_m(x_1, \ldots, x_t)$ be real polynomials, each of degree at most $d$. Then there are at most $(4edm/t)^t$ consistent sign–assignments of the form (14).

We consider the class of quantum circuits of size $N$ with $d$–bit gates computing $n$–variable Boolean functions. Without loss of generality, we can assume that $n'$, the number of input wires of such circuits, is at most $d \cdot N$. We define an equivalence relation $\simeq$ on such circuits: we write $C_1 \simeq C_2$ if and only if $C_1$ and $C_2$ differ only in the label of their gates; in another word, $C_1$ and $C_2$ have the same underlying graph but the corresponding gates in these circuits may compute different unitary operations. The number of different equivalence classes is at most $$(n')^d N \leq (dN)^{dN} = 2^{O(N \log N)}.$$

Now we find an upper bound for the number of different Boolean functions that can be computed by circuits in the same equivalence class. Fix an equivalence class $\mathcal{E}$. We use the variables $a_1 + ib_1, a_2 + ib_2, \ldots, a_\mu + ib_\mu$, where $\mu = d^2 N$, to denote the entries of the matrices of the gates of a circuit $C$ in $\mathcal{E}$. By
substituting appropriate values to the variables $a_1, \ldots, a_\mu, b_1, \ldots, b_\mu$, we get all circuits in $E$. On input $\alpha = (\alpha_1, \ldots, \alpha_n) \in \{0, 1\}^n$, the probability that $C$ outputs 1 can be represented by a real polynomial

$$P_\alpha(a_1, \ldots, a_\mu, b_1, \ldots, b_\mu).$$

The degree of $P_\alpha$ is at most $2N$. There are $2^n$ polynomials $P_\alpha$ and the number of different Boolean functions can be computed by $C$ by changing the unitary operators of its gates is at most the number of different consistent sign–assignments to the following system:

$$P_\alpha(a_1, \ldots, a_\mu, b_1, \ldots, b_\mu) - \frac{2}{3},$$
$$P_\alpha(a_1, \ldots, a_\mu, b_1, \ldots, b_\mu) - \frac{1}{3},$$

for $\alpha \in \{0, 1\}^n$. By Theorem 7.2, this number is bounded from the above by

$$\left(\frac{4e(2N)^{2^n+1}}{2\mu}\right)^{2\mu} = 2^{O(nN) + O(N \log N)}.$$

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