# Quantum data and control made easier 

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#### Abstract

In this paper we define nQML , a functional quantum programming language that follows the "quantum data and control" paradigm. In comparison to Altenkirch and Grattage's QML, the control constructs of nQML are simpler and can implement quantum algorithms more directly and naturally. We avoid the unnecessary complexities of a linear type system by using types that carry the address of qubits in the quantum state. We provide a denotational semantics over density matrices and unitary transformations, inspired by Selinger's semantics for QPL. Our semantics leads naturally to an interpreter for nQML, written in Haskell.


Keywords: Functional quantum programming language, type system, denotational semantics.

## 1 Introduction

In the years following the discovery of Shor's factoring algorithm [11] and Grover's algorithm for database search [5] the field of quantum computations has attracted much scientific interest. Unlike classical algorithms, quantum algorithms are almost invariably studied at a low level, involving quantum circuits and their properties. The fact that reasoning about quantum circuits is no easier than reasoning about their classical counterparts has given rise to quantum programming languages, that is, languages that allow programmers to implement quantum algorithms and make use of the added power of the quantum computational model, while respecting its special restrictions. In this paper we present such a language named nQML.

Our main focus in the design of $n Q M L$ is to give programmers sufficient expressive power to implement quantum algorithms easily, while preventing them from

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breaking the rules of quantum computation. nQML is a high-level functional language based on the concept of "quantum data and control". It includes constructs which allow any unitary transformation to be expressed as a program in nQML quite naturally, more or less using the same notation that is used by the designers of quantum algorithms. It also permits quantum measurements to be carried out at any point during the execution of a program.

The relative ease of use comes at the cost of putting aside a number of important practical issues, such as the existence of imperfect quantum hardware, the need for quantum error correction and the fact that every quantum program will eventually have to be implemented as a quantum circuit using only a finite set of quantum gates and, therefore, some of the unitary transformations that nQML allows will have to be approximated. Similar problems were a source of concern for the founders of the classical programming model many decades ago. Fortunately they have been resolved and their solutions have been abstracted in such a way that people who use modern high-level programming languages do not need to know anything about them. We believe that the same can and must be done for the quantum programming languages of the future and adopt the approach that such issues should be tackled not by the designer and users of a quantum programming language, but by the architect of a quantum computer, the designer of its operating system and, to a lesser extent, the designer of the compiler.
nQML admits a simple type system and denotational semantics. By simple, we mean that both use structures and techniques that are typical in the study of classical programming languages of similar size and complexity. They should therefore be easily accessible to readers with a basic knowledge of programming language semantics and an elementary understanding of the quantum computation model. The main novelty of nQML's type system is that the type of a quantum expression conveys information which reveals the exact qubits of the quantum state in which the expression's value resides. Qubit aliasing is allowed in such a way that the "no cloning" and "no dropping" principles are not violated. Programmers have the look-and-feel of a classical programming language, without linearity restrictions.

The denotational semantics of nQML is based on the use of density matrices to describe quantum states. The meaning of a well-typed nQML program is a function from density matrices to density matrices and describes the program's effect on an arbitrary quantum input state. Well-typed programs which conduct no measurements ${ }^{5}$ are also assigned a meaning in the form of a unitary matrix which describes the transformation they perform on the quantum state. The execution of a nQML program can be seen as a sequence of steps which affect the quantum state either by allocating new qubits, by applying unitary transformations to existing qubits or by measuring existing qubits. Our semantics leads to a straightforward implementation for nQML in the form of an interpreter written in Haskell. ${ }^{6}$ The interpreter, quite obviously, simulates quantum computations in a classical computer and takes exponential time.

The rest of the paper is structured as follows. Section 2 discusses related work.

[^1]In Section 3 we describe the syntax and semantics of nQML. Section 4 contains a number of examples, while Section 5 concludes with our final remarks. The appendix contains the complete formal definition of $n$ QML.

## 2 Related work

Starting with Knill's conventions for quantum pseudocode [6], several quantum programming languages have been proposed and an excellent survey of the emerging field can be found in [2]. Among the most notable are Ömer's QCL, an imperative language with quantum primitives and automatic quantum scratch space management [7], and Sanders and Zuliani's qGCL, an extension of Dijkstra's guarded command language [8]. Moreover, van Tonder has proposed a $\lambda$-calculus for higherorder quantum programs without measurements [12]. It is not clear however how this calculus corresponds to lower-level descriptions of quantum computations, such as quantum circuits.

Selinger's QPL is a language following the paradigm "quantum data, classical control" [9]. It is functional in nature, although from a programmer's point of view it looks more imperative than functional. QPL allows the programmer to access both classical and quantum memory and includes high-level features such as loops and recursion. Program control in QPL is strictly classical and quantum branching can only be implemented indirectly with appropriate unitary transformations. The denotational semantics of QPL is given in the form of superoperators on density matrices. A higher-order extension of QPL in the form of a quantum lambda calculus has also been proposed by Selinger and Valiron [10].

On the other hand, Altenkirch and Grattage's QML is a functional language that follows the paradigm "quantum data and control" $[1,3,4]$. QML comes with a linear type system which prohibits implicit weakening, which would lead to implicit measurements and quantum collapse. Variables in QML correspond to wires in the produced quantum circuit and thus have to be shared implicitly when they are used in several places in a program so as not to break the "no cloning" rule. The sharing of wires is also monitored by the linear type system. The semantics of QML assigns to every well-typed program a quantum circuit. QML's if ${ }^{\circ}$ operator implements the notion of quantum control and is the only available means of performing unitary transformation. The two branches of an if ${ }^{\circ}$ must be "orthogonal" quantum expressions, in order to preserve the reversibility of pure quantum computations.

The nature of our $n Q M L$ is inspired from QML, the main addition being the quantum transformation construct $|e\rangle \rightarrow x, x^{\prime} . c$ which will be described in Section 3. Its type system, although not linear, is an adaptation of Altenkirch and Grattage's type system. The semantics of nQML, however, is very much in the spirit of Selinger's denotational semantics for QPL.

## 3 The language nQML

The complete syntax of $n Q M L$ is given in the following grammar. It is assumed that $x$ is a variable identifier and $\lambda$ is a complex constant. The grammar defines two syntactic classes. Quantum expressions are denoted by $e$; they represent quantum
programs and their syntax is similar to that of QML. Classical expressions are denoted by $c$; they are only needed in the quantum transformation construct $|e\rangle \rightarrow$ $x, x^{\prime} . c$ and they can represent two types of information: a structure of classical bits or a complex number.

```
\(e::=x \mid\left\{(\lambda)\right.\) qfalse \(+\left(\lambda^{\prime}\right)\) qtrue \(\} \mid\) let \(x=e_{1}\) in \(e_{2}\)
    \(\left|\quad\left(e_{1}, e_{2}\right)\right|\) let \(\left(x_{1}, x_{2}\right)=e_{1}\) in \(e_{2}\)
    \(\mid \quad\) if \(e\) then \(e_{1}\) else \(e_{2} \mid\) ifm \(e\) then \(e_{1}\) else \(\left.e_{2}| | e\right\rangle \rightarrow x, x^{\prime} \cdot c\)
\(c::=x|\lambda|\) let \(x=c_{1}\) in \(c_{2}\left|\left(c_{1}, c_{2}\right)\right|\) let \(\left(x_{1}, x_{2}\right)=c_{1}\) in \(c_{2} \mid\) int \(c\)
    \(\left|c_{1}+c_{2}\right| c_{1}-c_{2}\left|c_{1} * c_{2}\right| c_{1} / c_{2}\left|c_{1}^{c_{2}}\right|\) if \(c\) then \(c_{1}\) else \(c_{2}\)
```

Variables in nQML are viewed as references to quantum information that is stored in a global quantum state. There are two types of quantum information: qubits and products. A new qubit is allocated in the quantum state when the superposition operator $\left\{(\lambda)\right.$ qfalse $+\left(\lambda^{\prime}\right)$ qtrue $\}$ is used, in the same way that new objects are allocated on the heap when a data constructor is used in a functional programming language. Products are introduced and eliminated with the constructs ( $e_{1}, e_{2}$ ) and let $\left(x_{1}, x_{2}\right)=e_{1}$ in $e_{2}$. nQML also features three control constructs:

- $\operatorname{ifm} e$ then $e_{1}$ else $e_{2}$ : It conducts a measurement on $e$, which must be of type qubit. Depending on the result, it executes one of its branches. It is similar to a classical random branching, based on a toss of a biased coin with probabilities depending on the state of the qubit being measured.
- if $e$ then $e_{1}$ else $e_{2}$ : It allows the programmer to perform quantum branching. If $e$, which must be of type qubit, is in a classical state, then the effect is what we would expect from ifm. But if $e$ is in a quantum superposition, the program proceeds in a quantum superposition of both branches, most likely creating entanglement among the qubits of the quantum state.
- $|e\rangle \rightarrow x, x^{\prime} . c:$ A generic means of expressing any unitary transformation, which has to be relied upon when a transformation can not be easily broken down to a series of controlled operations, expressible with if. Its advantage is that, rather than forcing programmers to precompute and provide the whole unitary matrix of the transformation, whose size is exponential in the number of qubits that it affects, it allows them to express that matrix as a complex function of the input and output state of the transformed qubits. This leads to a succinct and clear expression of many useful quantum algorithms, such as the quantum Fourier transform that is described in Section 4.

In quantum pseudocode notation, all unitary transformations can be expressed in the form:

$$
|i\rangle \rightarrow \sum_{j=0}^{2^{n}-1} f(i, j)|j\rangle
$$

where $f(i, j)$ is a function of the input state $i$ of the quantum register and its output state $j$. The construct $|e\rangle \rightarrow x, x^{\prime} . c$ allows the programmers to use precisely this natural notation: the classical variables $x$ and $x^{\prime}$ denote the register's input and output state and the classical expression $c$ denotes the function's body.

From this notation, if the function $f$ is known, the unitary matrix can be easily constructed by taking $S_{j, i}=f(i, j)$. Of course, not all functions $f$ result
in unitary matrices and the type system of nQML cannot decide whether the resulting transformation is indeed unitary. The type system of Altenkirch and Grattage's QML is able to do that, at the expense of making the size of the program exponential and complicating the typing with orthogonality constraints.

### 3.1 The type system of nQML

There are two kinds of types: quantum types $(\tau)$ and classical types $(\phi)$. For each quantum expression, the type system of nQML keeps track of the exact qubits of the state in which the value of this expression is stored. This information is stored in the types. It is used to make sure that the same qubit cannot be used twice in a transformation, thus allowing qubit aliasing without breaking the "no cloning" rule.

$$
\begin{aligned}
\tau & ::=\operatorname{qbit}[n] \mid \tau_{1} \otimes \tau_{2} \\
\phi & ::=\operatorname{bit}\left|\phi_{1} \times \phi_{2}\right| \text { complex }
\end{aligned}
$$

For example, an expression has type qbit[5] if its value is stored in the 5th qubit of the state.

For each quantum type $\tau$, we define $\mathcal{C}(\tau)$ to be the corresponding classical type; no quantum types correspond to complex. We denote by $|\mathcal{C}(\tau)|$ the size, in classical bits, of the classical type corresponding to $\tau$ and by $q \operatorname{bits}(\tau)$ the set of the state's qubits that are used by expressions of type $\tau$. For example, qbits $(\mathbf{q b i t}[4] \otimes$ $\operatorname{qbit}[2])=\{2,4\}$. A quantum type $\tau$ is called pure if its representation uses distinct qubits. Notice that, in general, $|q \operatorname{bits}(\tau)| \leq|\mathcal{C}(\tau)|$, the two being equal if and only if the type $\tau$ is pure. A quantum type environment $\Gamma$ is a mapping of variables to quantum types and, similarly, a classical type environment $\Delta$ is a mapping of variables to classical types. $\left.\Gamma\right|_{k}$ denotes the environment $\Gamma$ restricted in such a way that it does not contain variables whose types use the state's $k$-th qubit.

The typing relation for nQML is denoted by $\Gamma ; n \vdash^{\alpha} e: \tau ; m$. More precisely, as in the type system of Altenkirch and Grattage's QML, there are two typing relations: one for pure quantum expressions (i.e. without measurements), denoted by $\Gamma ; n \vdash^{\circ}$ $e: \tau ; m$, and one for arbitrary quantum expressions, denoted by $\Gamma ; n \vdash e: \tau ; m$. We refer to both by allowing the superscript ${ }^{\alpha}$ to be either ${ }^{\circ}$ or empty. As the types of nQML convey information regarding the position of qubits in the quantum state, the typing relation is forced to process and propagate such information. In $\Gamma ; n \vdash^{\alpha}$ $e: \tau ; m$, the natural number $n$ appearing on the left side of the relation stands for the number of qubits of the original quantum state, before $e$ starts evaluating. Obviously, for all pairs $\left(x: \tau_{x}\right) \in \Gamma$ it must be $q b i t s\left(\tau_{x}\right) \subseteq\{0, \ldots n-1\}$. The natural number $m$ appearing on the right side of the relation stands for the number of new qubits, that are allocated during the evaluation of $e$. The final quantum state after $e$ has been evaluated has $n+m$ qubits and, obviously again, it must be $q b i t s(\tau) \subseteq\{0, \ldots n+m-1\}$.

The typing rules for nQML follow Altenkirch and Grattage's QML, with the exception that the type system is not linear and qubit information must be processed. For example, the typing rule for quantum superposition plans for the allocation of one new qubit and uses its position in the returned type.

$$
\frac{|\lambda|^{2}+\left|\lambda^{\prime}\right|^{2}=1}{\Gamma ; n \vdash^{\circ}\left\{(\lambda) \text { qfalse }+\left(\lambda^{\prime}\right) \text { qtrue }\right\}: \text { qbit }[n] ; 1}
$$

Rules with more than one quantum expression must carefully combine the newly allocated qubits, e.g.

$$
\frac{\Gamma ; n \vdash^{\alpha} e_{1}: \tau_{1} ; m_{1} \quad \Gamma ; n+m_{1} \vdash^{\alpha} e_{2}: \tau_{2} ; m_{2}}{\Gamma ; n \vdash^{\alpha}\left(e_{1}, e_{2}\right): \tau_{1} \otimes \tau_{2} ; m_{1}+m_{2}} \text { (PROD) }
$$

The most complex of nQML's typing rules are those for the control constructs. We explain two of them below. In a quantum branching expression if $e$ then $e_{1}$ else $e_{2}$, the control qubit must not be used in the two branches. This restriction is necessary to simplify the semantics of if and eliminate the need for orthogonal branches. Unitary transformations which cannot easily be described as quantum controlled operations have their own dedicated construct in nQML. Notice also that the number of newly allocated qubits takes the maximum of the two branches.

$$
\begin{gathered}
\Gamma ; n \vdash^{\alpha} e: \underset{\text { qbit }[k] ; m}{ } \frac{\left.\Gamma\right|_{k} ; n+m \vdash^{\circ} e_{1}: \tau ;\left.m_{1} \quad \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{2}: \tau ; m_{2}}{\Gamma ; n \vdash^{\alpha} \text { if } e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right)} \text { (IF) }
\end{gathered}
$$

The typing rule for nQML's new construct $|e\rangle \rightarrow x, x^{\prime} . c$ is also straightforward. A unitary transformation is performed on the quantum bits where the value of expression $e$ is stored. The type $\tau$ of this expression must be pure, to obey the "no cloning" rule. In the classical expression $c$ which determines the contents of the transformation, the two variables $x$ and $x^{\prime}$ are bound to the classical value of the expression. The type of both is $\mathcal{C}(\tau)$.

$$
\begin{equation*}
\frac{\Gamma ; n \vdash^{\alpha} e: \tau ; m \quad \text { pure }(\tau) \quad x: \mathcal{C}(\tau), x^{\prime}: \mathcal{C}(\tau) \vdash c: \text { complex }}{\Gamma ; n \vdash^{\alpha}|e\rangle \rightarrow x, x^{\prime} . c: \tau ; m} \tag{TRANS}
\end{equation*}
$$

The typing $\Delta \vdash c: \phi$ of classical expressions presents no difficulties.

### 3.2 The denotational semantics of $n Q M L$

Our denotational semantics for nQML uses density matrices for representing the quantum state. The semantic domain $\mathbf{S}(n) \subset \mathbb{C}^{2^{n} \times 2^{n}}$ contains density matrices. The meaning of an arbitrary well-typed expression $e$ with a type derivation $\Gamma ; n \vdash$ $e: \tau ; m$ is a function of type $\mathbf{S}(n) \rightarrow \mathbf{S}(n+m)$; it maps an input quantum state of $n$ qubits to an output quantum state of $n+m$ qubits. Pure quantum expressions that perform no measurements can be assigned unitary transformations as meanings. We denote by $\mathbf{T}(n) \subset \mathbb{C}^{2^{n} \times 2^{n}}$ the domain of unitary transformation matrices. If $e$ is a well-typed pure quantum expression with a type derivation $\Gamma ; n \vdash^{\circ} e: \tau ; m$, then its meaning is a unitary transformation matrix of type $\mathbf{T}(n+m)$. The semantics of embedding pure quantum expressions in impure quantum expressions is given below. The tensor product of $A: \mathbf{S}(n)$ with the matrix $\Delta_{m}$ appropriately expands the state with $m$ new qubits which are initialized with zeroes.

EMB:

$$
\begin{gathered}
\llbracket \Gamma ; n \vdash e: \tau ; m \rrbracket(A)=T\left(A \otimes \Delta_{m}\right) T^{*} \\
\text { where } T=\llbracket \Gamma ; n \vdash^{\circ} e: \tau ; m \rrbracket
\end{gathered}
$$

The use of a variable has no effect on the state, as variables are just references. However, superpositions extend the state by allocating a new qubit and appropriately
initializing it.
VAR: $\quad \llbracket \Gamma ; n \vdash^{\circ} x: \tau ; 0 \rrbracket=\mathbb{I}_{n}$
SUP:

$$
\begin{aligned}
& \llbracket \Gamma ; n \vdash^{\circ}\left\{(\lambda) \text { qfalse }+\left(\lambda^{\prime}\right) \text { qtrue }\right\}: \text { qbit }[n] ; 1 \rrbracket= \\
& \mathbb{I}_{n} \otimes\left(\begin{array}{cc}
\lambda & \lambda^{\prime} \\
\lambda^{\prime} & -\lambda
\end{array}\right)
\end{aligned}
$$

The semantics of the let construct, product introduction and elimination is straightforward and very similar. In each of them, evaluation begins with the evaluation of $e_{1}$ and continues with the evaluation of $e_{2}$ on the new state. The impure cases are very similar. ${ }^{7}$
$L E T^{\circ}: \quad \llbracket \Gamma ; n \vdash^{\circ}$ let $x=e_{1}$ in $e_{2}: \tau ; m_{1}+m_{2} \rrbracket=T_{2}\left(T_{1} \otimes \mathbb{I}_{m_{2}}\right)$

$$
\text { where } \begin{aligned}
& T_{1}=\llbracket \Gamma ; n \vdash^{\circ} e_{1}: \tau_{1} ; m_{1} \rrbracket \\
& \\
& T_{2}=\llbracket \Gamma, x: \tau_{1} ; n+m_{1} \vdash^{\circ} e_{2}: \tau ; m_{2} \rrbracket
\end{aligned}
$$

The case of if is slightly more complicated. Evaluation begins with the condition. The matrices that correspond to the two branches are calculated and their (inexistent) effect on the control bit is removed by using the auxiliary function except. Then, the two expressions are executed conditionally, with $e$ as the control qubit. The impure case is again very similar.

$$
\begin{aligned}
& I F^{\circ}: \quad \llbracket \Gamma ; n \vdash^{\circ} \text { if } e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right) \rrbracket= \\
& T_{c}\left(T \otimes \mathbb{I}_{\max }\left(m_{1}, m_{2}\right)\right) \\
& \text { where } T=\llbracket \Gamma ; n \vdash^{\circ} e: \mathbf{q b i t}[k] ; m \rrbracket \\
& T_{1}=\left.\llbracket \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{1}: \tau ; m_{1} \rrbracket \\
& T_{2}=\left.\llbracket \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{2}: \tau ; m_{2} \rrbracket \\
& T_{1}^{\prime}=\operatorname{except}\left(k, T_{1}\right) \otimes \mathbb{I}_{\max \left(m_{1}, m_{2}\right)-m_{1}} \\
& T_{2}^{\prime}=\operatorname{except}\left(k, T_{2}\right) \otimes \mathbb{I}_{\max \left(m_{1}, m_{2}\right)-m_{2}} \\
& T_{c}=\operatorname{cond}\left(k, T_{1}^{\prime}, T_{2}^{\prime}\right)
\end{aligned}
$$

Surprisingly, the measuring conditional ifm is more straightforward. The condition is evaluated and then the corresponding qubit is measured. The auxiliary function measure returns the two density matrices that correspond to collapsing a qubit to a classical state. Then the two branches are combined. Each branch is evaluated on the corresponding result state of the measurement and their sum is the total result.

$$
\text { IFM: } \quad \begin{array}{r}
\llbracket \Gamma ; n \vdash \operatorname{ifm} e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right) \rrbracket(A)= \\
B_{1} \otimes \Delta_{\max \left(m_{1}, m_{2}\right)-m_{1}}+B_{2} \otimes \Delta_{\max }\left(m_{1}, m_{2}\right)-m_{2} \\
\text { where } B=\llbracket \Gamma ; n \vdash e: \operatorname{qbit}[k] ; m \rrbracket(A) \\
\left(B_{t}, B_{f}\right)=\text { measure }(k, B) \\
B_{1}=\llbracket \Gamma ; n+m \vdash e_{1}: \tau ; m_{1} \rrbracket\left(B_{t} \otimes \Delta_{m_{1}}\right) \\
B_{2}=\llbracket \Gamma ; n+m \vdash e_{2}: \tau ; m_{2} \rrbracket\left(B_{f} \otimes \Delta_{m_{2}}\right)
\end{array}
$$

Finally, in the semantics of $|e\rangle \rightarrow x, x^{\prime} . c$ the described unitary transformation $C$ is computed. As $C$ only applies to the qubits used by $e$, it must be properly expanded to apply to the complete state.

[^2]```
TRANS \({ }^{\circ}: \quad \llbracket \Gamma ; n \vdash^{\circ}|e\rangle \rightarrow x, x^{\prime} . c: \tau ; m \rrbracket=T_{c} T\)
    where \(T_{c}=\operatorname{expand}(n, q b i t s(\tau), C)\)
    \(T=\llbracket \Gamma ; n \vdash^{\circ} e: \tau ; m \rrbracket\)
    \(C_{j, i}=\llbracket x: \mathcal{C}(\tau), x^{\prime}: \mathcal{C}(\tau) \vdash c:\) complex \(\rrbracket(\rho)\)
    where \(\rho=\rho_{0}\left\{x \mapsto \operatorname{val}_{\tau}(i)\right\}\left\{x^{\prime} \mapsto \operatorname{val}_{\tau}(j)\right\}\)
    for all \(0 \leq i, j<2^{k}\), where \(k=|q b i t s(\tau)|\)
```

Again, the semantics of classical expressions is standard and presents no difficulty.

## 4 Examples

We now present a few example nQML programs of varying complexity. We start with two useful operators in quantum programming: not and had, standing respectively for quantum negation and the Hadamard transformation. Both can be applied to any expression $q$ of type qbit $[n]$.

```
\(\operatorname{not}(q) \equiv|q\rangle \rightarrow x, x^{\prime}\).if \(x^{\prime}=x\) then 0 else 1
\(\operatorname{had}(q) \equiv|q\rangle \rightarrow x, x^{\prime}\). if \(x\) then (if \(x^{\prime}\) then \(-\frac{1}{\sqrt{2}}\) else \(\frac{1}{\sqrt{2}}\) ) else \(\frac{1}{\sqrt{2}}\)
```

These simple transformations may seem a bit awkward at first but now that we have defined them we can easily use them in conjunction with the quantum conditional construct to define more complex transformations. For example, controlled quantum negation of $p$ by $q$ can be defined as:

```
cnot(q,p) \equiv if q then not (p) else p
```

where $\operatorname{not}(p)$ is defined as above.
This leads us to our first nQML program: an implementation of Deutsch's algorithm. In this algorithm we are presented with a black box classical one-bit boolean function and we want to decide whether it is balanced, in which case we return 1 , or constant, in which case we return 0 . We assume that the unknown function is somehow included in our program and we write $f(q)$ for the application of that function to a quantum parameter $q$. By using the definition of had and not given above, we arrive to the following program.

```
\(\operatorname{deutsch}(f) \equiv \operatorname{let}(i, j)=\left(\left\{\left(\frac{1}{\sqrt{2}}\right)\right.\right.\) qfalse \(+\left(\frac{1}{\sqrt{2}}\right)\) qtrue \(\}\),
    \(\left\{\left(\frac{1}{\sqrt{2}}\right)\right.\) qfalse \(+\left(-\frac{1}{\sqrt{2}}\right)\) qtrue \(\left.\}\right)\) in
let \(r=\) if \(f(i)\) then \(\operatorname{not}(j)\) else \(j\) in
had(i)
```

The program's result is stored in variable $i$. This variable is used as the first operand of our branching operator, after $f$ is applied to it. When $f$ is a constant function and therefore $f(i)$ has a classical value, $i$ will be unaffected by the execution of the branching and its result after the Hadamard transform will be 0 . When, however, $f$ is balanced, i.e. it is the identity or the negation function, even though its application will have no direct effect on $i$, the use of $i$ as a control bit for $j$ 's negation means that the two variables interact non-classically.

Let us now see a few more examples that demonstrate the power of $|e\rangle \rightarrow x, x^{\prime} . c$. Addition of a constant to a $n$-bit quantum register modulo $2^{n}$, which is typically denoted by $|r\rangle \rightarrow|r+c\rangle$ in quantum pseudocode, can be implemented as:

$$
\operatorname{add}(r, c) \equiv|r\rangle \rightarrow x, x^{\prime} . \text { if } \operatorname{int} x^{\prime}=\operatorname{int} x+c \text { then } 1 \text { else } 0
$$

Any other permutation of base states can easily be implemented in a similar manner. The implementation of the quantum Fourier transform for $n$ qubits contained in register $r$ is:

```
fourier \((r, n) \equiv|r\rangle \rightarrow x, x^{\prime} \cdot 1 / 2^{n} * e^{2 * \pi * i * x * x^{\prime} / 2^{n}}\)
```

which is derived in a straightforward way from the transform's definition. In Selinger's QPL, one can do the same by applying the unitary matrix $S$ corresponding to the quantum Fourier transform to the quantum register $r$, using the construct $r *=S$. However, unless some sophisticated language is used in combination with QPL to represent unitary transformations, the programmer has to use a precalculated $S$ and, as its size is $2^{n} \times 2^{n}$, the size of the program increases exponentially. The same is true in the case of Altenkirch and Grattage's QML, where the transform can be implemented by a tree of height $n$ containing nested if $^{\circ}$ branches; the size of the program is again exponential in $n$.

As a last example, let us see an implementation of Grover's fast database search. Assuming that $c$ denotes the value we are searching for, we first need to implement the query and diffusion operators.

```
query \((q) \quad \equiv|q\rangle \rightarrow x, x^{\prime}\). if \(x=x^{\prime}\)
    then (if int \(x=c\) then -1 else 1)
    else 0
diffusion \((q, n) \equiv|q\rangle \rightarrow x, x^{\prime}\). if \(x=x^{\prime}\) then \(-1+2 / 2^{n}\) else \(2 / 2^{n}\)
```

Let us consider the most simple application of Grover's algorithm: searching in a space of size 4 ( $n=2$ qubits). Even though $O(\sqrt{n})$ applications of the two operators are generally needed to obtain high probability, in this special case one application is enough to produce the correct result with certainty:

```
grover \(\equiv\) let \(q_{1}=\left\{\left(\frac{1}{\sqrt{2}}\right)\right.\) qfalse \(+\left(\frac{1}{\sqrt{2}}\right)\) qtrue \(\}\) in
    let \(q_{2}=\left\{\left(\frac{1}{\sqrt{2}}\right)\right.\) qfalse \(+\left(\frac{1}{\sqrt{2}}\right)\) qtrue \(\}\) in
    let \(q s=\left(q_{1}, q_{2}\right)\) in
    diffusion(query (qs),2)
```

Assuming that the element we were looking for was $c=2$, the Haskell interpreter that implements our semantics produces the following state (density matrix) of two qubits:

$$
\left(\begin{array}{cccc}
0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0 \\
0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0 \\
0.0:+0.0 & 0.0:+0.0 & 0.9999999999999997:+0.0 & 0.0:+0.0 \\
0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0 & 0.0:+0.0
\end{array}\right)
$$

where $\alpha:+\beta$ is Haskell's notation for the complex number $\alpha+\beta i$. From it, we can easily verify that the correct answer was found: the register $q s$ is in the classical state $|10\rangle$, allowing for numerical errors.

## 5 Conclusion

Quantum programming is today more or less at the same point in its history as classical programming was in the 1940s. The hardware is non existent or faulty. The semantics of quantum programming languages is understood either at a very low level of abstraction, using quantum gates and circuits, or at a very high level of abstraction, using tensor products in categories of Hilbert spaces. One thing that is different, though, is our experience of more than half a century in the theory and practice of classical programming languages. It is this experience that must be put into work if, sometime in the future, quantum programming languages are going to be what classical programming languages are today. Quantum programming must exploit the advantages of the quantum computational model, putting aside its peculiarities and insignificant details, so that programmers can add two "quantum integers" and obtain another "quantum integer" without, for example, having to think about the reversibility of this computation.

It can be argued that our work takes the "quantum data and control" paradigm a very small step further towards simplicity. We have defined nQML, a new functional quantum programming language, inspired by Altenkirch and Grattage's QML and following the "quantum data and control" paradigm. The type system of nQML keeps track of the use of qubits in expressions and avoids the complexities of linear type systems. Its semantics is inspired by Selinger's semantics for QPL. It is a simple denotational semantics with density matrices and unitary transformations as the semantic domains, which leads naturally to a simple implementation, in the form of an interpreter written in Haskell. Furthermore, the $|e\rangle \rightarrow x, x^{\prime} . c$ construct allows quantum algorithms to be implemented in a more direct and natural way.

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## A Formal definition of nQML

## A. 1 Syntax

```
e ::=x | {(\lambda) qfalse + (\lambda')}\mathrm{ qtrue } | let }x=\mp@subsup{e}{1}{}\mathrm{ in }\mp@subsup{e}{2}{
    (e, e, e2) | let (x, 利)= e
    if e then e}\mp@subsup{e}{1}{}\mathrm{ else e}\mp@subsup{e}{2}{}|\mathrm{ ifm e then }\mp@subsup{e}{1}{}\mathrm{ else }\mp@subsup{e}{2}{}||e\rangle->x,\mp@subsup{x}{}{\prime}.
```




## A. 2 Typing

Types: quantum and classical
$\tau::=\mathbf{q b i t}[n] \mid \tau_{1} \otimes \tau_{2}$
$\phi::=$ bit $\left|\phi_{1} \times \phi_{2}\right|$ complex

From quantum to classical types
$\mathcal{C}(\mathbf{q b i t}[n])=$ bit
$\mathcal{C}\left(\tau_{1} \otimes \tau_{2}\right)=\mathcal{C}\left(\tau_{1}\right) \times \mathcal{C}\left(\tau_{2}\right)$

Size of classical types
$\mid \mathcal{C}($ qbit $[n]) \mid=1$
$\left|\mathcal{C}\left(\tau_{1} \otimes \tau_{2}\right)\right|=\left|\mathcal{C}\left(\tau_{1}\right)\right|+\left|\mathcal{C}\left(\tau_{2}\right)\right|$
$\mid$ complex $\mid=$ undefined

Qubits used by a quantum type
$q \operatorname{bits}(\tau) \quad: \quad \mathcal{P}(\mathbb{N})$
qbits $(\mathbf{q b i t}[n])=\{n\}$
$q \operatorname{bits}\left(\tau_{1} \otimes \tau_{2}\right) \quad=\quad q \operatorname{bits}\left(\tau_{1}\right) \cup \operatorname{qbits}\left(\tau_{2}\right)$

Pure quantum types

$$
\frac{\operatorname{pure}(\mathbf{q b i t}[n])}{\operatorname{pure}\left(\tau_{1}\right)} \quad \operatorname{pure}\left(\tau_{2}\right) \quad \operatorname{qbits}\left(\tau_{1}\right) \cap q b i t s\left(\tau_{2}\right)=\emptyset \quad 1
$$

Type environments: quantum and classical
$\Gamma \quad: \quad$ a finite set of pairs of the form $(x: \tau)$
$\Delta \quad: \quad$ a finite set of pairs of the form $(x: \phi)$
$\left.\Gamma\right|_{k}=\{(x: \tau) \in \Gamma \mid k \notin q b i t s(\tau)\}$

Typing relation for quantum expressions

$$
\Gamma ; n \vdash^{\alpha} e: \tau ; m
$$

where ${ }^{\alpha}$ is empty or ${ }^{\circ}$

$$
\begin{gather*}
\frac{\Gamma ; n \vdash^{\circ} e: \tau ; m}{\Gamma ; n \vdash e: \tau ; m}(E M B) \quad \frac{(x: \tau) \in \Gamma}{\Gamma ; n \vdash^{\circ} x: \tau ; 0}(V A R) \\
\frac{|\lambda|^{2}+\left|\lambda^{\prime}\right|^{2}=1}{\Gamma ; n \vdash^{\circ}\left\{(\lambda) \text { qfalse }+\left(\lambda^{\prime}\right) \text { qtrue }\right\}: \text { qbit }[n] ; 1} \quad(S U P)  \tag{SUP}\\
\frac{\Gamma ; n \vdash^{\alpha} e_{1}: \tau_{1} ; m_{1} \quad \Gamma, x: \tau_{1} ; n+m_{1} \vdash^{\alpha} e_{2}: \tau ; m_{2}}{\Gamma ; n \vdash^{\alpha} \text { let } x=e_{1} \operatorname{in} e_{2}: \tau ; m_{1}+m_{2}}(L E T) \\
\frac{\Gamma ; n \vdash^{\alpha} e_{1}: \tau_{1} ; m_{1} \quad \Gamma ; n+m_{1} \vdash^{\alpha} e_{2}: \tau_{2} ; m_{2}}{\Gamma ; n \vdash^{\alpha}\left(e_{1}, e_{2}\right): \tau_{1} \otimes \tau_{2} ; m_{1}+m_{2}}(P R O D)
\end{gather*}
$$

$$
\begin{gathered}
\Gamma ; n \vdash^{\alpha} e_{1}: \tau_{1} \otimes \tau_{2} ; m_{1} \quad \Gamma, x_{1}: \tau_{1}, x_{2}: \tau_{2} ; n+m_{1} \vdash^{\alpha} e_{2}: \tau ; m_{2} \\
\hline \Gamma ; n \vdash^{\alpha} \text { let }\left(x_{1}, x_{2}\right)=e_{1} \text { in } e_{2}: \tau ; m_{1}+m_{2} \\
\Gamma ; n \vdash^{\alpha} e: \text { qbit }[k] ; m \\
\frac{\left.\Gamma\right|_{k} ; n+m \vdash^{\circ} e_{1}: \tau ;\left.m_{1} \quad \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{2}: \tau ; m_{2}}{\Gamma ; n \vdash^{\alpha} \text { if } e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right)} \text { (IF) } \\
\frac{\Gamma ; n \vdash e: \operatorname{qbit}[k] ; m \quad \Gamma ; n+m \vdash e_{1}: \tau ; m_{1} \quad \Gamma ; n+m \vdash e_{2}: \tau ; m_{2}}{\Gamma ; n \vdash \operatorname{ifm} e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right)} \text { (IFM) } \\
\frac{\Gamma ; n \vdash^{\alpha} e: \tau ; m \quad p u r e(\tau) \quad x: \mathcal{C}(\tau), x^{\prime}: \mathcal{C}(\tau) \vdash c: \text { complex }}{\Gamma ; n \vdash^{\alpha}|e\rangle \rightarrow x, x^{\prime} . c: \tau ; m} \text { (TRANS) }
\end{gathered}
$$

Typing relation for classical expressions

## A. 3 Semantics

Semantic domains

| $\mathbf{S}(n)$ | $=\left\{A \in \mathbb{C}^{2} \times 2^{n} \mid A\right.$ is a density matrix $\}$ |
| :--- | :--- |
| $\mathbf{T}(n)$ | $=\left\{T \in \mathbb{C}^{2} \times 2^{n} \mid T\right.$ is unitary $\}$ |
| $\llbracket \Delta \rrbracket$ | $=\Pi x:$ Var. $\llbracket \Delta(x) \rrbracket$ |
| $\llbracket$ bit $\rrbracket$ | $=\mathbb{B}$ |
| $\llbracket \phi_{1} \times \phi_{2} \rrbracket$ | $=\llbracket \phi_{1} \rrbracket \times \llbracket \phi_{2} \rrbracket$ |
| $\llbracket$ complex $\rrbracket$ | $=\mathbb{C}$ |

Semantics of pure quantum expressions

$$
\llbracket \Gamma ; n \vdash^{\circ} e: \tau ; m \rrbracket: \mathbf{T}(n+m)
$$

VAR: $\quad \llbracket \Gamma ; n \vdash^{\circ} x: \tau ; 0 \rrbracket=\mathbb{I}_{n}$
SUP:
$L E T^{\circ}$ :

$$
\llbracket \Gamma ; n \vdash^{\circ}\left\{(\lambda) \text { qfalse }+\left(\lambda^{\prime}\right) \text { qtrue }\right\}: \mathbf{q b i t}[n] ; 1 \rrbracket=\mathbb{I}_{n} \otimes\left(\begin{array}{cc}
\lambda & \lambda^{\prime} \\
\lambda^{\prime} & -\lambda
\end{array}\right)
$$

$\llbracket \Gamma ; n \vdash^{\circ}$ let $x=e_{1}$ in $e_{2}: \tau ; m_{1}+m_{2} \rrbracket=T_{2}\left(T_{1} \otimes \mathbb{I}_{m_{2}}\right)$
where $T_{1}=\llbracket \Gamma ; n \vdash^{\circ} e_{1}: \tau_{1} ; m_{1} \rrbracket$
$\mathrm{PROD}{ }^{\circ}$. $\quad \Gamma ; n{ }_{2}=\llbracket \Gamma, x: \tau_{1} ; n+m_{1} \vdash^{\circ} e_{2}: \tau ; m_{2} \rrbracket$

$$
\llbracket \Gamma ; n \vdash^{\circ}\left(e_{1}, e_{2}\right): \tau_{1} \otimes \tau_{2} ; m_{1}+m_{2} \rrbracket=T_{2}\left(T_{1} \otimes \mathbb{I}_{m_{2}}\right)
$$

where $T_{1}=\llbracket \Gamma ; n \vdash^{\circ} e_{1}: \tau_{1} ; m_{1} \rrbracket$

$$
T_{2}=\llbracket \Gamma ; n+m_{1} \vdash^{\circ} e_{2}: \tau_{2} ; m_{2} \rrbracket
$$

$\operatorname{LETPROD}{ }^{\circ}: \quad \llbracket \Gamma ; n \vdash^{\circ}$ let $\left(x_{1}, x_{2}\right)=e_{1}$ in $e_{2}: \tau ; m_{1}+m_{2} \rrbracket=T_{2}\left(T_{1} \otimes \mathbb{I}_{m_{2}}\right)$
where $T_{1}=\llbracket \Gamma ; n \vdash^{\circ} e_{1}: \tau_{1} \otimes \tau_{2} ; m_{1} \rrbracket$
$T_{2}=\llbracket \Gamma, x_{1}: \tau_{1}, x_{2}: \tau_{2} ; n+m_{1} \vdash^{\circ} e_{2}: \tau ; m_{2} \rrbracket$
$I F^{\circ}:$

$$
\llbracket \Gamma ; n \vdash^{\circ} \text { if } e \text { then } e_{1} \text { else } e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right) \rrbracket=
$$

$$
T_{c}\left(T \otimes \mathbb{I}_{\max \left(m_{1}, m_{2}\right)}\right)
$$

where $T=\llbracket \Gamma ; n \vdash^{\circ} e: \mathbf{q b i t}[k] ; m \rrbracket$
$T_{1}=\left.\llbracket \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{1}: \tau ; m_{1} \rrbracket$
$T_{2}=\left.\llbracket \Gamma\right|_{k} ; n+m \vdash^{\circ} e_{2}: \tau ; m_{2} \rrbracket$
$T_{1}^{\prime}=\operatorname{except}\left(k, T_{1}\right) \otimes \mathbb{I}_{\max }\left(m_{1}, m_{2}\right)-m_{1}$
$T_{2}^{\prime}=\operatorname{except}\left(k, T_{2}\right) \otimes \mathbb{I}_{\max }\left(m_{1}, m_{2}\right)-m_{2}$
$T_{c}=\operatorname{cond}\left(k, T_{1}^{\prime}, T_{2}^{\prime}\right)$

$$
\begin{aligned}
& \frac{(x: \phi) \in \Delta}{\Delta \vdash x: \phi}(\text { var }) \quad \frac{}{\Delta \vdash \lambda: \text { complex }} \text { (const) } \\
& \frac{\Delta \vdash c_{1}: \phi_{1} \Delta, x: \phi_{1} \vdash c_{2}: \phi}{\Delta \vdash \operatorname{let} x=c_{1} \text { in } c_{2}: \phi} \text { (let) } \frac{\Delta \vdash c_{1}: \phi_{1} \Delta \vdash c_{2}: \phi_{2}}{\Delta \vdash\left(c_{1}, c_{2}\right): \phi_{1} \times \phi_{2}} \text { (prod) } \\
& \frac{\Delta \vdash c_{1}: \phi_{1} \times \phi_{2} \quad \Delta, x_{1}: \phi_{1}, x_{2}: \phi_{2} \vdash c_{2}: \phi}{\Delta \vdash \operatorname{let}\left(x_{1}, x_{2}\right)=c_{1} \text { in } c_{2}: \phi} \text { (letprod) } \\
& \frac{\Delta \vdash c: \mathcal{C}(\tau)}{\Delta \vdash \operatorname{int} c: \text { complex }} \text { (int) } \frac{\Delta \vdash c: \operatorname{bit} \Delta \vdash c_{1}: \phi \quad \Delta \vdash c_{2}: \phi}{\Delta \vdash \operatorname{if} c \text { then } c_{1} \text { else } c_{2}: \phi} \text { (if) } \\
& \frac{\Delta \vdash c_{1}: \text { complex } \quad \Delta \vdash c_{2}: \text { complex } \quad o p \in\left\{+,-, *, /,^{\wedge}\right\}}{\Delta \vdash c_{1} \text { op } c_{2}: \text { complex }} \text { (op) }
\end{aligned}
$$

TRANS ${ }^{\circ}$ :
$\llbracket \Gamma ; n \vdash^{\circ}|e\rangle \rightarrow x, x^{\prime} . c: \tau ; m \rrbracket=T_{c} T$
where $\quad T_{c}=\operatorname{expand}(n, \operatorname{qbits}(\tau), C)$

$$
T=\llbracket \Gamma ; n \vdash^{\circ} e: \tau ; m \rrbracket
$$

$C_{j, i}=\llbracket x: \mathcal{C}(\tau), x^{\prime}: \mathcal{C}(\tau) \vdash c:$ complex $\rrbracket(\rho)$
where $\rho=\rho_{0}\left\{x \mapsto \operatorname{val}_{\tau}(i)\right\}\left\{x^{\prime} \mapsto \operatorname{val}_{\tau}(j)\right\}$
for all $0 \leq i, j<2^{k}$, where $k=|q \operatorname{bits}(\tau)|$

Semantics of impure quantum expressions

$$
\llbracket \Gamma ; n \vdash e: \tau ; m \rrbracket: \mathbf{S}(n) \rightarrow \mathbf{S}(n+m)
$$

| EMB: | $\begin{gathered} \llbracket \Gamma ; n \vdash e: \tau ; m \rrbracket(A)=T\left(A \otimes \Delta_{m}\right) T^{*} \\ \text { where } T=\llbracket \Gamma ; n \vdash^{\circ} e: \tau ; m \rrbracket \end{gathered}$ |
| :---: | :---: |
| LET : | $\llbracket \Gamma ; n \vdash$ let $x=e_{1}$ in $e_{2}: \tau ; m_{1}+m_{2} \rrbracket(A)=B_{2}$ where $\quad B_{1}=\llbracket \Gamma ; n \vdash e_{1}: \tau_{1} ; m_{1} \rrbracket(A)$ |
| PROD : | $\begin{gathered} B_{2}=\llbracket \Gamma, x: \tau_{1} ; n+m_{1} \vdash e_{2}: \tau ; m_{2} \rrbracket\left(B_{1}\right) \\ \llbracket \Gamma ; n \vdash\left(e_{1}, e_{2}\right): \tau_{1} \otimes \tau_{2} ; m_{1}+m_{2} \rrbracket(A)=B_{2} \\ \text { where } B_{1}=\llbracket \Gamma ; n \vdash e_{1}: \tau_{1} ; m_{1} \rrbracket(A) \end{gathered}$ |
| LETPROD : | $\begin{aligned} B_{2} & =\llbracket \Gamma ; n+m_{1} \vdash e_{2}: \tau_{2} ; m_{2} \rrbracket\left(B_{1}\right) \\ \llbracket \Gamma ; n \vdash \text { let }\left(x_{1}, x_{2}\right) & =e_{1} \text { in } e_{2}: \tau ; m_{1}+m_{2} \rrbracket(A)= \\ \text { where } B_{1} & =\llbracket \Gamma ; n \vdash e_{1}: \tau_{1} \otimes \tau_{2} ; m_{1} \rrbracket(A) \\ B_{2} & =\llbracket \Gamma, x_{1}: \tau_{1}, x_{2}: \tau_{2} ; n+m_{1} \vdash e_{2}: \tau ; m_{2} \rrbracket\left(B_{1}\right) \end{aligned}$ |
| IF : |  |
| IFM: | $\llbracket \Gamma ; n \vdash \operatorname{ifm} e$ then $e_{1}$ else $e_{2}: \tau ; m+\max \left(m_{1}, m_{2}\right) \rrbracket(A)=$ $B_{1} \otimes \Delta_{\max \left(m_{1}, m_{2}\right)-m_{1}}+B_{2} \otimes \Delta_{\max \left(m_{1}, m_{2}\right)-m_{2}}$ where $B=\llbracket \Gamma ; n \vdash e: \mathbf{q b i t}[k] ; m \rrbracket(A)$ $\left(B_{t}, B_{f}\right)=\text { measure }(k, B)$ $B_{1}=\llbracket \Gamma ; n+m \vdash e_{1}: \tau ; m_{1} \rrbracket\left(B_{t} \otimes \Delta_{m_{1}}\right)$ <br> $B_{2}=\llbracket \Gamma ; n+m \vdash e_{2}: \tau ; m_{2} \rrbracket\left(B_{f} \otimes \Delta_{m_{2}}\right)$ |
| TRANS : | $\begin{aligned} \llbracket \Gamma ; n \vdash\|e\rangle \rightarrow & x, x^{\prime} \cdot c: \tau ; m \rrbracket(A)=T_{c} B T_{c}^{*} \\ \text { where } & T_{c}=\operatorname{expand}(n, q \operatorname{bits}(\tau), C) \\ & B=\llbracket \Gamma ; n \vdash e: \tau ; m \rrbracket(A) \\ & C_{j, i}=\llbracket x: \mathcal{C}(\tau), x^{\prime}: \mathcal{C}(\tau) \vdash c: \operatorname{complex} \rrbracket(\rho) \\ & \text { where } \rho=\rho_{0}\left\{x \mapsto v a l_{\tau}(i)\right\}\left\{x^{\prime} \mapsto \operatorname{val}(j)\right\} \\ & \text { for all } 0 \leq i, j<2^{k}, \text { where } k=\|q b i t s(\tau)\| \end{aligned}$ |

Semantics of classical expressions

$$
\llbracket \Delta \vdash c: \phi \rrbracket: \llbracket \Delta \rrbracket \rightarrow \llbracket \phi \rrbracket
$$

```
var: \(\quad \llbracket \Delta \vdash x: \phi \rrbracket(\rho)=\rho(x)\)
const: \(\quad \llbracket \Delta \vdash \lambda:\) complex \(\rrbracket(\rho)=\lambda\)
let: \(\quad \llbracket \Delta \vdash\) let \(x=c_{1}\) in \(c_{2}: \phi \rrbracket(\rho)=\llbracket \Delta, x: \phi_{1} \vdash c_{2}: \phi \rrbracket\left(\rho^{\prime}\right)\)
    where \(\rho^{\prime}=\rho\left\{x \mapsto \llbracket \Delta \vdash c_{1}: \phi_{1} \rrbracket(\rho)\right\}\)
prod: \(\quad \llbracket \Delta \vdash\left(c_{1}, c_{2}\right): \phi_{1} \times \phi_{2} \rrbracket(\rho)=\left(\llbracket \Delta \vdash c_{1}: \phi_{1} \rrbracket(\rho), \llbracket \Delta \vdash c_{2}: \phi_{2} \rrbracket(\rho)\right)\)
letprod: \(\quad \llbracket \Delta \vdash \operatorname{let}\left(x_{1}, x_{2}\right)=c_{1}\) in \(c_{2}: \phi \rrbracket(\rho)=\)
        \(\llbracket \Delta, x_{1}: \phi_{1}, x_{2}: \phi_{2} \vdash c_{2}: \phi \rrbracket\left(\rho^{\prime}\right)\)
    where \(\left(v_{1}, v_{2}\right)=\llbracket \Delta \vdash c_{1}: \phi_{1} \times \phi_{2} \rrbracket(\rho)\)
    \(\rho^{\prime}=\rho\left\{x \mapsto v_{1}\right\}\left\{y \mapsto v_{2}\right\}\)
int: \(\quad \llbracket \Delta \vdash \operatorname{int} c:\) complex \(\rrbracket(\rho)=\operatorname{code}_{\tau}(\llbracket \Delta \vdash c: \mathcal{C}(\tau) \rrbracket(\rho))\)
op: \(\quad \llbracket \Delta \vdash c_{1}\) op \(c_{2}\) : complex \(\rrbracket(\rho)=\)
    \(\llbracket \Delta \vdash c_{1}\) : complex \(\rrbracket(\rho)\) op \(\llbracket \Delta \vdash c_{2}\) : complex \(\rrbracket(\rho)\)
if:
\(\llbracket \Delta \vdash\) if \(c\) then \(c_{1}\) else \(c_{2}: \phi \rrbracket(\rho)=\)
    \(\left\{\begin{array}{l}\llbracket \Delta \vdash c_{1}: \phi \rrbracket(\rho), \text { if } \llbracket \Delta \vdash c: \operatorname{bit} \rrbracket(\rho)=\text { true } \\ \llbracket \Delta \vdash c_{2}: \phi \rrbracket(\rho), \text { if } \llbracket \Delta \vdash c: \operatorname{bit} \rrbracket(\rho)=\text { false }\end{array}\right.\)
```

Auxiliary functions
$\mathbb{I}_{n} \quad: \quad$ the identity matrix of size $2^{n} \times 2^{n}$
$\Delta_{n} \quad$ : a matrix of size $2^{n} \times 2^{n}$ with all zeroes and a 1 in the top-left corner
except $: \quad \mathbb{N} \times \mathbf{S}(n+1) \rightarrow \mathbf{S}(n)$

```
\(\operatorname{except}\left(0,\left(\begin{array}{c|c}A & \mathbb{O} \\ \hline \mathbb{O} & A\end{array}\right)=\quad=\quad A\right.\)
\(\operatorname{except}\left(k+1,\left(\begin{array}{c|c}A & B \\ \hline C & D\end{array}\right)=\left(\frac{\operatorname{except}(k, A) \mid \operatorname{except}(k, B)}{\operatorname{except}(k, C) \mid \operatorname{except}(k, D)}\right)\right.\)
cond : \(\quad \mathbb{N} \times \mathbf{S}(n) \times \mathbf{S}(n) \rightarrow \mathbf{S}(n+1)\)
\(\operatorname{cond}(0, T, F)=\left(\begin{array}{c|c}F & \mathbb{O} \\ \hline \mathbb{O} \mid T\end{array}\right)\)
\(\operatorname{cond}\left(k+1,\left(\begin{array}{c|c}T_{A} & T_{B} \\ \hline T_{C} & T_{D}\end{array}\right),\left(\begin{array}{l|l}F_{A} & F_{B} \\ \hline F_{C} & F_{D}\end{array}\right)\right)=\left(\begin{array}{c}\operatorname{cond}\left(k, T_{A}, F_{A}\right) \mid \operatorname{cond}\left(k, T_{B}, F_{B}\right) \\ \hline \operatorname{cond}\left(k, T_{C}, F_{C}\right) \\ \operatorname{cond}\left(k, T_{D}, F_{D}\right)\end{array}\right)\)
measure : \(\mathbb{N} \times \mathbf{S}(n+1) \rightarrow \mathbf{S}(n+1) \times \mathbf{S}(n+1)\)
measure \(\left.\left(0,\left(\begin{array}{l|l|l}A & \mathbb{O} \\ \hline \mathbb{O} & D\end{array}\right)\right)=\left(\begin{array}{c|c|c}\mathbb{O} & \mathbb{O} \\ \hline \mathbb{O} & D\end{array}\right),\left(\begin{array}{l|l}A & \mathbb{O} \\ \hline \mathbb{O} & \mathbb{O}\end{array}\right)\right)\)
measure \(\left(k+1,\left(\begin{array}{c|c}A & \mathbb{O} \\ \hline \mathbb{O} & A\end{array}\right)=\left(\left(\begin{array}{l|l|}T_{A} & T_{B} \\ \hline T_{C} & T_{D}\end{array}\right),\left(\begin{array}{l|l}F_{A} & F_{B} \\ \hline F_{C} & F_{D}\end{array}\right)\right)\right.\)
    where \(\left(T_{A}, F_{A}\right)=\) measure \((k, A)\)
        \(\left(T_{B}, F_{B}\right)=\) measure \((k, B)\)
        \(\left(T_{C}, F_{C}\right)=\) measure \((k, C)\)
        \(\left(T_{D}, F_{D}\right)=\) measure \((k, D)\)
expand \(\quad: \quad \Pi n: \mathbb{N} . \Pi S: \mathcal{P}(\mathbb{N}) . \mathbf{T}(|S|) \rightarrow \mathbf{T}(n)\)
\(\operatorname{expand}(n, S, T)=\operatorname{expa}_{0}(n, S, T)\)
    where \(\operatorname{expa}_{n}(n, S, T)=T\)
        \(\operatorname{expa}_{k}\left(n, S,\left(\begin{array}{c|c}A & B \\ \hline C \mid D\end{array}\right)=\left(\frac{\operatorname{expa}_{k+1}(n, S, A) \mid \operatorname{expa}}{k+1}(n, S, C),\left(\operatorname{expa} a_{k+1}(n, S, B) \mid \operatorname{expa} a_{k+1}(n, S, D) \quad\right)\right.\right.\)
                if \(k<n\) and \(k \in S\)
            \(\operatorname{expa}_{k}(n, S, T)=\mathbb{I}_{1} \otimes \operatorname{expa}_{k+1}(n, S, T)\)
                if \(k<n\) and \(k \notin S\)
code \(_{\tau} \quad: \quad \llbracket \mathcal{C}(\tau) \rrbracket \rightarrow \mathbb{N}\)
\(\operatorname{code}_{\mathbf{q b i t}[k]}(b)=\left\{\begin{array}{l}1, \text { if } b=\text { true } \\ 0, \text { if } b=\text { false }\end{array}\right.\)
\(\operatorname{code}_{\tau_{1} \otimes \tau_{2}}\left(v_{1}, v_{2}\right)=2^{k} \operatorname{code}_{\tau_{1}}\left(v_{1}\right)+\operatorname{code}_{\tau_{2}}\left(v_{2}\right)\)
    where \(k=\left|\mathcal{C}\left(\tau_{2}\right)\right|\)
val \(_{\tau} \quad: \quad \mathbb{N} \rightarrow \llbracket \mathcal{C}(\tau) \rrbracket\)
\(\operatorname{val}_{\text {qbit }[k]}(n)=\left\{\begin{array}{l}\text { true }, \text { if } n=1 \\ \text { false, if } n=0\end{array}\right.\)
\(\operatorname{val}_{\tau_{1} \otimes \tau_{2}}(n)=\left(\operatorname{val}_{\tau_{1}}\left(n / 2^{k}\right), \operatorname{val}_{\tau_{2}}\left(n \bmod 2^{k}\right)\right)\)
    where \(k=\left|\mathcal{C}\left(\tau_{2}\right)\right|\)
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[^1]:    5 In the sequel, such programs will be called "pure" quantum programs, for short.
    6 The source code of the interpreter is available from ftp://ftp.softlab.ntua.gr/pub/users/nickie/ papers/nqml06.code.tar.gz.

[^2]:    7 It can easily be proved that the semantics of pure and impure quantum expressions is consistent with the embedding rule. For example, the meaning is the same if EMB is applied separately to two pure expressions and then PROD is applied to the result, or if EMB is applied once to the result of PROD.

