Workshop on Quantum Computation using IBM Q, Part I

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- Our First Quantum Circuit!
- A deeper look into quantum algorithms





My warmest thanks to (1/5)

Our sponsors - Thank you very much!







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My warmest thanks to (3/5)

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Fronteras de la Ciencia and SNI grants







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Access to the D-Wave II quantum annealer @ NASA Ames Center







My warmest thanks to (5/5)

My warmest thanks to you all for taking part in this initiative!







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Our goal

To contribute towards making quantum computing

- i) A solid research field in Mexico and all Latin America
 - ii) An engine of regional economic development





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Who we are

QIP Group at Tecnológico de Monterrey (1/2)

- We founded the field of quantum computation in Mexico.
- Quantum algorithms.
- Quantum Walks.
- Quantum annealing.
- Emerging fields of quantum computation, e.g. Quantum Image Processing and Quantum Machine Learning.
- Classical/digital simulation of quantum algorithms .
- Novel methods for training quantum scientists and quantum engineers.





Who we are

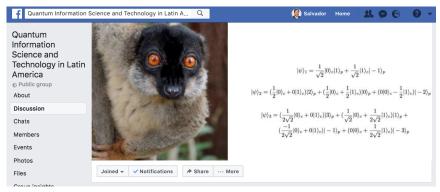
QIP Group at Tecnológico de Monterrey (2/2)

- Marco Lanzagorta. NRL, US Navy, USA. Quantum radar, quantum walks, quantum image processing, quantum annealing algorithms.
- Catherine McGeoch. D-Wave Systems, Canada. Quantum annealing algorithms.
- William de la Cruz de los Santos. UAEM. Quantum annealing algorithms, quantum computer vision.
- Mauricio García. IBM. Simulating and running quantum algorithms on IBM Q platforms.
- Enrique Hernández-Lemus. INMEGEN. Quantum algorithms for De novo Genome Sequence Assembly.
- Fei Yan, Changchun University, China and Ahmed Abdel-Latif, Menoufia University, Egypt. Quantum image processing.





Who we are



Join our FB Group!

https://www.facebook.com/groups/152996132018274/







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What is quantum computation?

Quantum computation, one of the most recent joint ventures between physics and computer science, may be defined as follows:

Quantum Computation is a multidisciplinary field focused on the development of computers and algorithms, *i.e. hardware* and *software*, based on the quantum mechanical properties of Nature.





QC, a multidisciplinary and pervasive field

- Quantum computation is transitioning from an emerging branch of science into a mature research field. A growing number of quantum scientists & engineers are devoting their efforts to identifying and developing cross-fertilising initiatives in fields such as:
 - Artificial intelligence
 - Machine learning
 - Computational geometry
 - Image processing





Many new players and interests

- Quantum computation is also an emerging market of advanced technology that is attracting:
 - Investors. Breakoff Capital (www.breakoffcapital.com), Quantum Valley Investments (http://quantumvalleyinvestments.com/).
 - Global Companies. Microsoft (https://www.stationq.com/), Google (https://plus.google.com/+QuantumAlLab), IBM (http://www.research.ibm.com/quantum/)
 - Entrepreneurs. 1 Qbit (http://1qbit.com/), Rigetti (http://rigetti.com/), D-Wave (http://www.dwavesys.com/).
 - Government. EU (https://ec.europa.eu/digital-single-market/en/news/european-commission-will-launch-eu1-billion-quantum-technologies-flagship), UK (http://uknqt.epsrc.ac.uk/), USA (https://www.whitehouse.gov/blog/2016/07/26/realizing-potential-quantum-information-science-and-advancing-high-performance).





Timeline (Theory)

Timeline (Theory)



Feynman QM computers



Deutsch Universal QTM



Shor Integer factorization



Grover Search an unsorted DB



Nishimori Quantum Annealing



DiVincenzo Criteria to construct a QC



Quantum Walks



Farhi Quantum Adiabatic Computation





Companies (1/2)

Companies (global)



Quantum Annealers (Optimization problems). SaS and direct selling. Classical simulation of quantum annealers

STATION



Topological quantum computation Classical simulators of quantum algorithms (Visual Studio-QC)



16, 25, 50 qubit universal computer. SaS. -Universal quantum computer Software platform for classical simulation of quantum - Quantum simulation of quantum systems algorithms and development environment.







Companies (2/2)

Companies (startups)



Quantum algorithms Machine learning



Quantum algorithms Quantum hardware



Unconventional Computing
Quantum algorithms







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Quantum algorithms: fast and furious! (1/4)

A goal of paramount importance in quantum computation is to determine **how** to manipulate quantum mechanical systems in order to develop algorithms that are *faster* than their classical counterparts.

Here, faster has two meanings:





Quantum algorithms: fast and furious! (2/4)

 Computational speed-up as a heritage of physical properties of quantum systems.

Example: quantum parallelism, i.e. quantum interference as a computational resource.

$$|0\rangle^{\otimes n}|0\rangle \overset{\hat{H}^{\otimes n}\hat{I}}{\longmapsto} \frac{1}{\sqrt{2^n}} \sum_{i=0}^{2^n-1} |i\rangle|0\rangle \overset{\hat{U}_f}{\longmapsto} \frac{1}{\sqrt{2^n}} \sum_{i=0}^{2^n-1} |i\rangle|f(i)\rangle$$





Quantum algorithms: fast and furious! (3/4)

 Development of algorithmic strategies based on quantum mechanical properties.

Example in classical computer science: for solving the TSP we may use a brute force algorithm O(n!) or the Bellman–Held–Karp algorithm $O(n^22^n)$.

Some preliminary results in quantum computation but much is yet to be done!





Quantum algorithms: fast and furious! (4/4)

Building quantum algorithms is a most challenging task because:

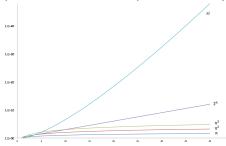
- a) First of all, we need a complex interesting problem (e.g., NP-complete, NP-hard).
- b) Quantum mechanics is a counterintuitive theory and intuition plays a key role in algorithm design.
- c) A competitive quantum algorithm must not only produce the solutions it is expected to, it also has to be more efficient, at least for some input values, than any classical algorithm (at least more efficient than existing classical algorithms).





Complex interesting problem?

Easy vs Difficult Algorithms (1 step/ns $\Rightarrow 3.15 \times 10^{22}$ steps/one million years)



Decision problems, P, NP, NP - complete, combinatorial optimisation problems, NP - hard.





Bits and Qubits - a quick introduction (1/6)

In order to motivate our discussion, let us briefly introduce the basic units of information in classical and quantum computation.







Bits and Qubits - a quick introduction (2/6)

In classical computation, information is stored and manipulated in the form of **bits**.

The mathematical structure of a classical bit is rather simple: it suffices to define two logical values, traditionally labelled as $\{0,1\}$ (i.e. a classical bit lives in a scalar space) and to relate those values to two different and mutually exclusive outcomes of a classical measurement.

Examples:



A relay and a transistor (Source: Wikimedia Commons)





Bits and Qubits - a quick introduction (3/6)

In quantum computation, information is stored, manipulated and measured in the form of **qubits** (= **quantum bit**).

A qubit is a vector that represents the state of two-state quantummechanical physical entity.



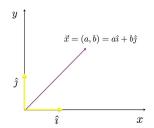


Bits and Qubits - a quick introduction (4/6)

Remember:

- $-\mathbb{R}^2$, the two-dimensional space studied in elementary school. Elements of \mathbb{R}^2 , denoted by \vec{x} , are vectors.
- Any vector $\vec{x}=(a,b)$ can be written as a linear combination of a basis, e.g.

$$\vec{x} = (a, b) = \vec{x} = (a, b) = a\hat{\imath} + b\hat{\jmath}$$









Bits and Qubits - a quick introduction (5/6)

In quantum computation, we use the following notation for denoting vectors:

$$\vec{x} = |x\rangle$$

So,

$$\vec{x} = a\hat{\imath} + b\hat{\jmath} \iff |x\rangle = a|i\rangle + b|j\rangle$$





Bits and Qubits - a quick introduction (6/6)

We shall shortly learn more about qubits but, before we do so, let us briefly focus on some key aspects of theoretical computer science.







But, before we go on...

Python, QSKIT, Jupiter: installing and troubleshooting.







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What is the ultimate objective of Theoretical Computer Science?

Theoretical Computer Science, traditionally thought of as a branch of mathematics, is divided into three fields linked by the question

What are the fundamental capabilities and limits of computers?







Fields of the Theory of Computation

<u>Automata theory.</u> Definitions and properties of abstract models of computation.

Computability theory. What problems can(not) be solved by computers?

Complexity theory. What makes some problems computationally hard and others easy?





Automata (1/3)

Automata are mathematical models of physical devices used to compute.

The definition of any automaton requires three components: states, language and transition rules.





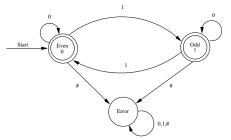
Automata (2/3)

- States are abstractions of the different stages an automaton would go through while computing (e.g., on and off states of a switch).
- A language is a set of strings made by concatenating characters from a given alphabet. For example, one could define a language from alphabet {a,b}, consisting of all strings that begin with 'a' and end with 'b'. This language would include 'ab', 'aabb', and 'ababab', but not 'ba' or 'bbb'.
- Finally, transition rules describe the internal dynamics of an automaton (i.e., how and when an automaton changes its state).





Automata (3/3) - An example of an automaton

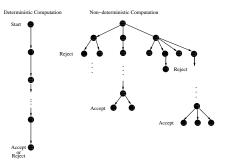


A finite automaton for parity computation. Double-circled states are accept states and single-circled state is a reject state. Input strings for $\mathcal R$ are any combination of characters taken from $\Sigma=\{0,1,\#\}$. Since an empty set of characters has no 1s, we set the initial parity of a string as even (this is why the state 'Even' has an arrow labeled as 'Start').





(Non)Deterministic Computation (1/3)



Deterministic automata follow a simple rule: given an input datum, every state of the computation is followed by only one state.







(Non)Deterministic Computation (2/3)

- Each and every automaton has deterministic and nondeterministic versions.
- Deterministic automata follow a simple rule : given an input datum, every state of the computation is followed by only one state.
- In contrast, when computing on a nondeterministic machine, a state could be followed by *many* states, depending on the input datum and the transition rules. Nondeterministic computation can be thought of as parallel computation with unlimited resources.





(Non)Deterministic Computation (3/3)

- Resources are always limited, hence nondeterministic computation may seem to be an unreasonable model.
- However, there are two good reasons to use this model: i) to determine if a problem is computable even if limitless resources are available, and ii) to define a key set of computable problems: NP problems
- An approximate realization of non-deterministic computation: probabilistic algorithms.





Turing Machines (1/3)

So far, Turing machines constitute the most accurate model of general purpose computers.



Turing Machines (drawing by Tom Dunne, American Scientist, March-April 2002)

The 'hardware' elements of a Deterministic Turing Machine (DTM) are a limitless memory tape (the tape is divided into squares or cells), and a scanner which consists of a read-write head *plus* a finite state control system. The scanner has two purposes: to read and write information on the cells of the tape as well as to control the state of the DTM.





Turing Machines (2/3)

There are Deterministic and Nondeterministic Turing machines.

DTMs are constrained to having access to only one computational state at a time, NTMs have access to as many states as needed.

Both Turing machines models are equally powerful in terms of computability.





Turing Machines (3/3)

Three important properties of DTMs:

- 1) There is a Turing Machine, known as **The Universal Turing Machine** (UTM), that can simulate the behavior of any other Turing machine.
- 2) The UTM and the von Neumann computer (our silicon-based computers, like laptops) are equivalent in terms of computability power. In other words, any problem that can be solved using a von Neumann computer can also be solved using the UTM.
- 3) Although highly nontrivial, an arbitrary program written in any computer language can be translated into a series of elementary steps in a DTM.





Algorithm Analysis

Algorithm analysis is divided into two parts:

- Algorithm Correctness.
- Algorithm resource consumption.







Algorithm Correctness

An algorithm A is correct if:

- a) [Design step] A successfully does what it is designed to do. This step usually involves using solid mathematical statements (e.g., theorems).
- b) [Implementation step] Implementation of A is robust and free of bugs.

Of course, easier said than done!





Algorithm analysis also requires thinking about the amount of resources that an algorithm will consume during its execution. More precisely,

We need to understand/estimate how resource requirements will **scale** as input size increases.





By Complexity Measure we mean a definition of efficiency that is platform-independent, instance-independent and of predictive value with respect to increasing input sizes.







A measure based on specific implementation details like computer hardware instruction sets (e.g., Intel or AMD processors) or the primitives of a particular computer language (e.g., Python or Lisp) would produce, for one and the same algorithm, different output values for different computer platforms.





Consequently, we will formulate abstract descriptions to outline algorithm behavior.

The key concept here is, as you may expect, the notion of **ele-mentary step**.





We shall analyze algorithm performance in two different scenarios:

- Worst case complexity. We are interested in estimating an upper bound on the maximum number of elementary steps that an algorithm must carry out. This is particularly important for pathological inputs.
- Average case complexity. We are interested in estimating the expected number of elementary steps that an algorithm must typically carry out. This is useful to estimate how an algorithm should perform 'in practice'.

In both cases, complexity if a function of input size |x|.





Worst-case complexity

Def. Worst-case complexity. Let D_n be the set of inputs of size n for the problem under consideration and $I \in D_n$.

Also, let t(I) be the number of elementary steps performed by the algorithm on input I. We define the function W by

$$W = \max\{t(I)|I \in D_n\}$$

W(n) is the maximum number of basic operations performed by the algorithm on any input of size n.





Average-case complexity

Def. Average-case complexity Let D_n be the set of inputs of size n for the problem under consideration and $I \in D_n$.

Also, let t(I) be the number of elementary steps performed by the algorithm on input I and p(I) be the probability that input I occurs. We define the function A by

$$A = \sum_{I \in D_n} p(I)t(I)$$





Problem. Let L be an array containing n different entries. If $x \in L$ return the index of x in L (that is, the array slot number in which x is contained). Return 0 otherwise (i.e., if $x \notin L$). Elementary step. Comparison of x with a list entry.

Algorithm

```
\begin{array}{l} index := 1 \\ \text{while } (index \leq n \text{ and } L[index] \neq x) \text{ do} \\ \mid index := index + 1 \\ \text{end} \\ \text{if } index > n \text{ then} \\ \mid index := 0 \\ \text{end} \end{array}
```





Worst case. Worst cases occur when x appears only in the last position in the list and when x is not in the list at all.

In both cases, we need to compare \boldsymbol{x} with all elements of L, i.e. execute \boldsymbol{n} comparisons. Hence,

$$W(n) = n$$





Average case. We shall analyze two cases:

- a) x is an element of L .
- b) x may not be an element of L.







- a) $x \in L$.
 - Suppose that x is equally likely to be in any particular position.
 - Let I_i , $i \in \{1, 2, \dots n\}$ represent the case where x appears in the i^{th} position of L.
 - Let $t(I_i)$ be the number of comparisons done (that is, the number of times the condition $L[index] \neq x$ is tested).
 - Note that, for $i \in \{1, 2, \dots n\}, t(I_i) = i$.

Thus,

$$A(n) = \sum_{i=1}^{n} p(I_i)t(I_i) = \sum_{i=1}^{n} \frac{1}{n}i = \frac{1}{n}\sum_{i=1}^{n} i = \frac{1}{n}\frac{n(n+1)}{2} = \frac{(n+1)}{2}$$

So,



$$A(n) = \frac{(n+1)}{2}$$



- b) x may not be an element of L.
 - Let α be the probability that $x \in L$ and $\beta = 1 \alpha$ the probability that $x \notin L$.
 - Suppose that if $x \in L$ then x is equally likely to be in any particular position.
 - Let I_i , $i \in \{1, 2, ..., n\}$ represent the case where x is in the i^{th} position of L.
 - Let I_{n+1} represent the case in which $x \notin L$.
 - Let $t(I_i)$ be the number of comparisons done.
 - Note that, for $i \in \{1, 2, \dots n\}, t(I_i) = i$ and $t(I_{n+1}) = n$.
 - Finally, for $i \in \{1, 2, \dots n\}, p(I_i) = \frac{\alpha}{n}$ and $p(I_{n+1}) = \beta$.





Thus,

$$A(n) = \sum_{i=1}^{n+1} p(I_i)t(I_i)$$

$$= \sum_{i=1}^{n} \frac{\alpha}{n}i + \beta I_{n+1}$$

$$= \frac{\alpha}{n} \sum_{i=1}^{n} i + \beta I_{n+1}$$

$$= \frac{\alpha}{n} \frac{n(n+1)}{2} + \beta I_{n+1}$$

$$= \frac{\alpha(n+1)}{2} + \beta I_{n+1}$$

So,



$$A(n) = \frac{\alpha(n+1)}{2} + \beta I_{n+1}$$



Intuitive Intro to Computational Complexity (1/10)

In complexity theory, for a problem $\ensuremath{\mathfrak{P}}$ that can be computed in a Turing machine, we are interested in answering the following question:

What is the minimum amount of time or energy / number of elementary steps that must be invested in order to solve \mathfrak{P} ?





Intuitive Intro to Computational Complexity (2/10)

Wait! What do we mean by elementary step?

- Machine language
- Assembly language
- High level language
- Abstract algorithm (like computing the inner product between two vectors).

The key point: polynomial transformations from one level of algorithm description to another.

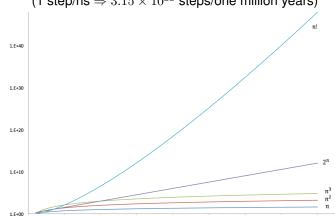




Intuitive Intro to Computational Complexity (3/10)

Easy vs Difficult Algorithms

(1 step/ns $\Rightarrow 3.15 \times 10^{22}$ steps/one million years)







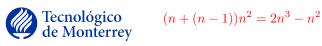


Intuitive Intro to Computational Complexity (4/10)

Example of an easy problem: square matrix multiplication. Assume addition and multiplication as elementary steps.

$$\begin{pmatrix} a_{11} & \cdots & a_{1n} \\ a_{21} & \cdots & a_{2n} \\ \vdots & & \vdots \\ a_{n1} & \cdots & a_{nn} \end{pmatrix} \begin{pmatrix} b_{11} & \cdots & b_{1n} \\ b_{21} & \cdots & b_{2n} \\ \vdots & & \vdots \\ b_{n1} & \cdots & b_{nn} \end{pmatrix} = \begin{pmatrix} \sum_{i=1}^{n} a_{1i}b_{i1} & \cdots & \sum_{i=1}^{n} a_{1i}b_{in} \\ \sum_{i=1}^{n} a_{2i}b_{i1} & \cdots & \sum_{i=1}^{n} a_{2i}b_{in} \\ \vdots & & \vdots \\ \sum_{n=1}^{n} a_{ni}b_{i1} & \cdots & \sum_{i=1}^{n} a_{ni}b_{in} \end{pmatrix}$$

The textbook multiplication of two $n \times n$ matrices requires n multiplications and n-1 additions per entry. Since the resulting matrix has n^2 entries, then the total amount of elementary steps required to multiply two $n \times n$ matrices is



$$(n + (n-1))n^2 = 2n^3 - n$$



Intuitive Intro to Computational Complexity (5/10)

For example, suppose that n=1000 and that a computer running the above-mentioned algorithm computes an elementary step in $1~\mu s$ (which is a rather slow computer). So,

$$2n^3 - n^2|_{n=10^3} = 2 \times 10^9 - 10^6 = 1999 \times 10^6$$
 elementary steps

Now, if each elementary step takes $1~\mu s$ then the total running time will be equal to

$$(1999\times 10^6 {\rm _{elementary\ steps}})(1\times 10^{-6}\frac{s}{{\rm _{elementary\ steps}}})=1999s=33.31\ min$$





Intuitive Intro to Computational Complexity (6/10)

In contrast, let us now analyze a key problem in theoretical computer science: the K-SAT problem.





Intuitive Intro to Computational Complexity (7/10)

K-SAT, a Fundamental NP-Complete Problem

Let $B = \{x_1, x_2, \dots, x_n, \bar{x_1}, \bar{x_2}, \dots \bar{x_n}\}$ be a set of Boolean variables. Also, let C_i be a disjunction of k elements of k and k be a conjunction of k clauses k

Question: Is there an assignment of Boolean variables in F that satisfies all clauses simultaneously, i.e. F=1?

For example:





Intuitive Intro to Computational Complexity (8/10)

Suppose $B=\{x_1,x_2,x_3,x_4,x_5,x_6,\bar{x_1},\bar{x_2},\bar{x_3},\bar{x_4},\bar{x_5},\bar{x_6}\}$ and

$$\begin{array}{lll} P & = & (\bar{x_1} \vee \bar{x_4} \vee \bar{x_5}) \wedge (\bar{x_2} \vee \bar{x_3} \vee \bar{x_4}) \wedge (x_1 \vee x_2 \vee \bar{x_5}) \wedge (x_3 \vee x_4 \vee x_5) \wedge \\ & (x_4 \vee x_5 \vee \bar{x_6}) \wedge (\bar{x_1} \vee \bar{x_3} \vee \bar{x_5}) \wedge (x_1 \vee \bar{x_2} \vee \bar{x_5}) \wedge (x_2 \vee \bar{x_3} \vee \bar{x_6}) \wedge \\ & (\bar{x_1} \vee \bar{x_2} \vee \bar{x_6}) \wedge (x_3 \vee \bar{x_5} \vee \bar{x_6}) \wedge (\bar{x_1} \vee \bar{x_2} \vee \bar{x_4}) \wedge (x_2 \vee x_3 \vee \bar{x_4}) \wedge \\ & (x_2 \vee x_5 \vee \bar{x_6}) \wedge (x_2 \vee \bar{x_3} \vee \bar{x_5}) \wedge (\bar{x_2} \vee \bar{x_3} \vee \bar{x_4}) \wedge (x_2 \vee x_3 \vee x_6) \wedge \\ & (\bar{x_1} \vee \bar{x_2} \vee \bar{x_3}) \wedge (\bar{x_1} \vee \bar{x_4} \vee \bar{x_5}) \wedge (\bar{x_3} \vee \bar{x_4} \vee \bar{x_6}) \wedge (\bar{x_4} \vee \bar{x_5} \vee x_6) \wedge \\ & (\bar{x_2} \vee x_3 \vee \bar{x_6}) \wedge (x_2 \vee x_5 \vee x_6) \wedge (x_3 \vee x_5 \vee \bar{x_6}) \wedge (\bar{x_1} \vee x_3 \vee \bar{x_6}) \wedge \\ & (x_3 \vee \bar{x_5} \vee x_6) \wedge (x_4 \vee x_5 \vee x_6) \wedge (x_1 \vee x_2 \vee \bar{x_3}) \end{array}$$

Finding the solutions (if any) of even a modest SAT instance can become difficult quite easily. In fact, only $\{x_1=0,x_2=1,x_3=0,x_4=1,x_5=0,x_6=0\}$ satisfies P.





Intuitive Intro to Computational Complexity (9/10)

Given an arbitrary instance F, we can always determine whether there is an assignment of Boolean variables such that F = 1.

To do so, we just have to substitute each and every possible value of x_i , $i \in \{1, \dots n\}$ in F. However, this procedure would require us to produce 2^n different assignments, i.e. an *exponential* number of different assignments.





Intuitive Intro to Computational Complexity (10/10)

For example, suppose that n=1000 (remember, n is the number of binary variables) and that a computer running the abovementioned brute force algorithm produces an assignment in 1fs, i.e. $10^{-15}s$ (a fast computer, indeed!)

Since $2^{1000}\approx 1.0715086072\times 10^{301}$ then the total amount of time to be invested in running our brute force algorithm would be about

$$1.0715086072 \times 10^{301} \times 10^{-15} s = 1.0715086072 \times 10^{286} s$$



 $3.39 \times 10^{278} \text{ years!}$



P, NP

Decision problems, a key concept in complexity theory, can be informally defined as yes/no questions.

The notion of decision problem is key to define the complexity classes known as P, NP and NP - complete. This is because decision problems can be put in correspondence with languages, being the latter a formal mathematical concept in computer science.





P, NP

– Class P. For each decision problem $\mathcal{A} \in \mathsf{P}$ there is a DTM M and a polynomial p(n) such that processing input ω of length n in M would be done in p(n) steps at most.





– Class NP. For each decision problem $\mathfrak{B}\in \mathsf{NP}$ there is an NTM N and a polynomial p(n) such that processing input ω of length n in N would be done in p(n) steps at most.





Theorem. If decision problem A is in NP then there is a polynomial p(n), where n is the size of an arbitrary input data, such that A can be solved by a deterministic algorithm having time complexity $O(2^{p(n)})$.





Equivalent definition for NP. A problem $\mathfrak B$ is in NP if there is a polynomial time algorithm designed to test whether a proposed solution to $\mathfrak B$ is correct.

Note: worrying about how hard it might be to find the solution is not part of this definition \odot

So, if only we could guess the right solution, we could then quickly test it.





It is well known that $P \subseteq NP$ and it is unknown whether $NP \subseteq P$.

Hence, a key research area in contemporary computer science is the study of decision problems in NP for which we only know inefficient algorithms, i.e. algorithms with time complexity functions that are not upper-bounded by polynomials.





Example 1 - NP problems /1/2)

Solving hard instances of the K-SAT problem

K-SAT

Let $B = \{x_1, x_2, \dots, x_n, \bar{x_1}, \bar{x_2}, \dots \bar{x_n}\}$ be a set of Boolean variables. Also, let C_i be a disjunction of k elements of k and k be a conjunction of k clauses k

Question: Is there an assignment of Boolean variables in F that satisfies all clauses simultaneously, i.e. F=1?





Example 1 - NP problems (2/2)

Suppose $B=\{x_1,x_2,x_3,x_4,x_5,x_6,\bar{x_1},\bar{x_2},\bar{x_3},\bar{x_4},\bar{x_5},\bar{x_6}\}$ and

$$\begin{array}{ll} P & = & (\bar{x_1} \vee \bar{x_4} \vee \bar{x_5}) \wedge (\bar{x_2} \vee \bar{x_3} \vee \bar{x_4}) \wedge (x_1 \vee x_2 \vee \bar{x_5}) \wedge (x_3 \vee x_4 \vee x_5) \wedge \\ & (x_4 \vee x_5 \vee \bar{x_6}) \wedge (\bar{x_1} \vee \bar{x_3} \vee \bar{x_5}) \wedge (x_1 \vee \bar{x_2} \vee \bar{x_5}) \wedge (x_2 \vee \bar{x_3} \vee \bar{x_6}) \wedge \\ & (\bar{x_1} \vee \bar{x_2} \vee \bar{x_6}) \wedge (x_3 \vee \bar{x_5} \vee \bar{x_6}) \wedge (\bar{x_1} \vee \bar{x_2} \vee \bar{x_4}) \wedge (x_2 \vee x_3 \vee \bar{x_4}) \wedge \\ & (x_2 \vee x_5 \vee \bar{x_6}) \wedge (x_2 \vee \bar{x_3} \vee \bar{x_5}) \wedge (\bar{x_2} \vee \bar{x_3} \vee \bar{x_4}) \wedge (x_2 \vee x_3 \vee x_6) \wedge \\ & (\bar{x_1} \vee \bar{x_2} \vee \bar{x_3}) \wedge (\bar{x_1} \vee \bar{x_4} \vee \bar{x_5}) \wedge (\bar{x_3} \vee \bar{x_4} \vee \bar{x_6}) \wedge (\bar{x_4} \vee \bar{x_5} \vee x_6) \wedge \\ & (\bar{x_2} \vee x_3 \vee \bar{x_6}) \wedge (x_2 \vee x_5 \vee x_6) \wedge (x_3 \vee x_5 \vee \bar{x_6}) \wedge (\bar{x_1} \vee x_3 \vee \bar{x_6}) \wedge \\ & (x_3 \vee \bar{x_5} \vee x_6) \wedge (x_4 \vee x_5 \vee x_6) \wedge (x_1 \vee x_2 \vee \bar{x_3}) \end{array}$$

Finding the solutions (if any) of even a modest SAT instance can become difficult quite easily. In fact, only $\{x_1=0,x_2=1,x_3=0,x_4=1,x_5=0,x_6=0\}$ satisfies P.

Notice that, for an binary string S, it would be possible to quickly determine whether S is a solution to an instance of the K-SAT problem. Tecnológico



Example 2 - Optimisation and NP problems (1/2)

Def. The Traveling Salesman Problem (optimization version)

INSTANCE: A finite set $C = \{c_1, c_2, \dots c_m\}$ of "cities" and a "distance" $d(c_i, c_j) \in \mathbb{N}$, the set of natural numbers.

QUESTION: Which is the shortest "tour" of all the cities in C, that is, an ordering $[c_{\Pi(1)},c_{\Pi(2)},\ldots,c_{\Pi(m)}]$ of C such that

$$\left[\sum_{i=1}^{m-1} d(c_{\Pi(i)}, c_{\Pi(i+1)})\right] + d(c_{\Pi(m)}, c_{\Pi(1)})$$

is minimum?

de Monterrey

Suppose we were given a tour T_i . Would it be possible to quickly check whether T_i is the shortest tour of all possible tours? Tecnológico



Example 2 - Optimisation and NP problems (2/2)

Def. The Traveling Salesman Problem (decision problem version)

INSTANCE: A finite set $C=\{c_1,c_2,\dots c_m\}$ of "cities" and a "distance" $d(c_i,c_j)\in\mathbb{N}$ and a bound $B\in\mathbb{N}$.

QUESTION: Is there a "tour" of all the cities in C having total length no more than B, that is, an ordering $[c_{\Pi(1)},c_{\Pi(2)},\ldots,c_{\Pi(m)}]$ of C such that $[\sum_{i=1}^{m-1}d(c_{\Pi(i)},c_{\Pi(i+1)})]+d(c_{\Pi(m)},c_{\Pi(1)})\leq B$?





Introductions to Computational Complexity for Physicists (1/2)

LIMITS OF COMPUTATION

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COMPUTATIONAL COMPLEXITY FOR PHYSICISTS

The theory of computational complexity has some interesting links to physics, in particular to quantum computing and statistical mechanics. This article contains an informal introduction to this theory and its links to physics.

S. Mertens. IEEE Journal on Computing in Science & Engineering vol. 4, no. 3, May/June 2002, pp 31-47







Introductions to Computational Complexity for Physicists (2/2)

CONTEMPORARY PHYSICS, 2018 https://doi.org/10.1080/00107514.2018.1450720





A cross-disciplinary introduction to quantum annealing-based algorithms

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ABSTRACT

A central goal in quantum computing is the development of quantum hardware and quantum algorithms in order to analyse challenging scientific and engineering problems. Research in quantum computation involves contributions from both physics and computer science; hence this article presents a concise introduction to basic concepts from both fields that are used in annealing-based quantum computation, an alternative to the more familiar quantum gate model. We introduce some concepts from computer science required to define difficult computational problems and to realise the potential relevance of quantum algorithms to find novel solutions to those problems. We introduce the structure of quantum annealing-based algorithms as well as two examples of this kind of algorithms for solving instances of the max-SAT and Minimum Multicut problems. An overview of the quantum annealing systems manufactured by D-Wave Systems is also presented.

ARTICLE HISTORY

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KEYWORDS

Quantum annealing; quantum algorithms; quantum computation







So far,

We have learned some basic notions of automata and complexity theory, certainly enough to realize the actual and potential relevance of quantum computation.

Let us now focus on the mathematics of quantum computation.





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Hilbert space

A **Hilbert space** \mathcal{H} is a (complete) complex inner-product vector space. An example of a Hilbert space is \mathbb{C}^n .





Kets and Bras

The Bra-Ket notation, also known as Dirac Notation, is a standard representation to describe quantum states.

The Dirac notation is widely used in quantum mechanics and quantum computation.

Let us now formally define the notions of Ket and Bra.





Kets and Bras

Kets. Let \mathcal{H} be a Hilbert space. A vector $\psi \in \mathcal{H}$ is denoted by $|\psi\rangle$ and it is referred to as a **ket**.

We can represent elements $|\psi\rangle$ of $\mathcal H$ as column vectors by choosing a basis for $\mathcal H$. For example, let $\mathcal H=\mathbb C^2$ and let us choose the vector basis $\{|0\rangle,|1\rangle\}$, where

$$|0\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix}$$
 and $|1\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$

Then, every element $|\psi\rangle\in\mathcal{H}$ can be written as

$$|\psi\rangle=\alpha|0\rangle+\beta|1\rangle=\alpha\begin{pmatrix}1\\0\end{pmatrix}+\beta\begin{pmatrix}0\\1\end{pmatrix}$$
, $\alpha,\beta\in\mathbb{C}$





Example of kets

$$|\psi\rangle=\left(rac{\frac{1}{\sqrt{2}}}{\frac{1}{\sqrt{2}}}
ight)\in\mathbb{C}^2$$
 may be written as

$$|\psi\rangle = \begin{pmatrix} \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{pmatrix} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 \\ 0 \end{pmatrix} + \frac{1}{\sqrt{2}} \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \frac{1}{\sqrt{2}} |0\rangle + \frac{1}{\sqrt{2}} |1\rangle$$





Exercise - Kets

Let
$$|+\rangle=\frac{|0\rangle+|1\rangle}{\sqrt{2}}$$
 and $|-\rangle=\frac{|0\rangle-|1\rangle}{\sqrt{2}}.$ Write $|\psi\rangle=\alpha|0\rangle+\beta|1\rangle$ in terms of $|+\rangle, |-\rangle.$





Answer to exercise - Kets

Let
$$|+\rangle = \frac{|0\rangle + |1\rangle}{\sqrt{2}}$$
 and $|-\rangle = \frac{|0\rangle - |1\rangle}{\sqrt{2}}$. Write $|\psi\rangle = \alpha|0\rangle + \beta|1\rangle$ in terms of $|+\rangle, |-\rangle$.

$$|+\rangle + |-\rangle = \frac{|0\rangle + |1\rangle}{\sqrt{2}} + \frac{|0\rangle - |1\rangle}{\sqrt{2}} = \frac{2|0\rangle}{\sqrt{2}} \Rightarrow |0\rangle = \frac{|+\rangle + |-\rangle}{\sqrt{2}}$$

Similarly,

$$|+\rangle - |-\rangle = \frac{|0\rangle + |1\rangle}{\sqrt{2}} - (\frac{|0\rangle - |1\rangle}{\sqrt{2}}) = \frac{2|1\rangle}{\sqrt{2}} \Rightarrow |1\rangle = \frac{|+\rangle - |-\rangle}{\sqrt{2}}$$

Hence,

$$|\psi\rangle=\alpha|0\rangle+\beta|1\rangle=\alpha^{\frac{|+\rangle+|-\rangle}{\sqrt{2}}}+\beta^{\frac{|+\rangle-|-\rangle}{\sqrt{2}}}=\frac{\alpha+\beta}{\sqrt{2}}|+\rangle+\frac{\alpha-\beta}{\sqrt{2}}|-\rangle$$

Therefore.



$$|\psi\rangle = \frac{\alpha+\beta}{\sqrt{2}}|+\rangle + \frac{\alpha-\beta}{\sqrt{2}}|-\rangle$$



Bras

Bras. Formally speaking, bras are functionals (i.e. functions of vector spaces into corresponding fields) and in practice, they can be thought of as row vectors:

$$|\psi\rangle = \alpha|0\rangle + \beta|1\rangle$$
 if and only if $|\psi\rangle = \alpha^*\langle 0| + \beta^*\langle 1|$





Bras

For example, let

$$|\psi\rangle = \frac{i}{\sqrt{2}} \begin{pmatrix} 1 \\ 0 \end{pmatrix} + \frac{1}{\sqrt{2}} \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} \frac{i}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{pmatrix} = \frac{i}{\sqrt{2}} |0\rangle + \frac{1}{\sqrt{2}} |1\rangle$$

The corresponding bra $\langle \psi |$ is

$$\langle \psi | = \frac{-i}{\sqrt{2}}(1, \ 0) + \frac{1}{\sqrt{2}}(0, \ 1) = (\frac{-i}{\sqrt{2}}, \frac{1}{\sqrt{2}}) = \frac{-i}{\sqrt{2}}\langle 0 | + \frac{1}{\sqrt{2}}\langle 1 |$$





Exercises - Bras

- 1. Compute $\langle +|$ and $\langle -|$
- 2. Let $|\psi\rangle=\alpha|0\rangle+\beta|1\rangle$ where $||\alpha||^2+||\beta||^2=1$. Does it follow that $\langle\psi|=\alpha^*\langle 0|+\beta^*\langle 1|$ where $||\alpha^*||^2+||\beta^*||^2=1$?





Answers to exercises - Bras

1. Compute $\langle +|$ and $\langle -|$

Answer:
$$\langle +|=\frac{\langle 0|+\langle 1|}{\sqrt{2}}$$
 and $\langle -|=\frac{\langle 0|-\langle 1|}{\sqrt{2}}$





Answers to exercises - Bras

2. Let $|\psi\rangle = \alpha|0\rangle + \beta|1\rangle$ where $||\alpha||^2 + ||\beta||^2 = 1$. Does it follow that $|\langle\psi| = \alpha^*\langle 0| + \beta^*\langle 1|$ where $||\alpha^*||^2 + ||\beta^*||^2 = 1$?

Answer:

- a) $|\psi\rangle = \alpha |0\rangle + \beta |1\rangle \Rightarrow \langle \psi| = \alpha^* \langle 0| + \beta^* \langle 1|$.
- b) Since $\alpha, \beta \in \mathbb{C}$ then let us write $\alpha = a + bi$ and $\beta = c + di$.
- c) Also, $\alpha^* = a bi$ and $\beta^* = c di$
- d) Furthermore, $||\alpha||^2=a^2+b^2$ and $||\beta||^2=c^2+d^2\Rightarrow$

$$||\alpha||^2 + ||\beta||^2 = a^2 + b^2 + c^2 + d^2 = 1$$

e) Finally, please note that $||\alpha^*||^2 = a^2 + b^2$ and $||\beta^*||^2 = c^2 + d^2 \Rightarrow ||\alpha^*||^2 + ||\beta^*||^2 = a^2 + b^2 + c^2 + d^2 = 1$

So, the answer is Yes, it does.





Summary of Kets and Bras

Thus, if ${\mathcal H}$ is an n-dimensional Hilbert space then

- A ket $|\psi\rangle\in\mathcal{H}$ can be represented as an n-dimensional column vector.
- Its corresponding bra $\langle \psi | \in \mathcal{H}^*$ can be seen as an n-dimensional row vector

 $|\psi\rangle\leftrightarrow\langle\psi|$ corresponds to transposition and conjugation.





Inner product

We can use the Dirac notation to make calculations. For example, $\langle \phi | \psi \rangle$ is the usual row-column matrix operator that computes the inner product in finite dimensional vector spaces.

For instance, let us take the representations of $|0\rangle$ and $|1\rangle$ given in previous slides

$$|0\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix} \text{ and } |1\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$$

Note that $|0\rangle \perp |1\rangle$ as well as the fact that both vectors have unitary norm. Consequently, the inner product of $|0\rangle$ and $|1\rangle$ must be zero and the inner product of each vector with itself must be equal to one:





Inner product

$$\langle \mathbf{0} | \mathbf{1} \rangle = (1, 0) \begin{pmatrix} 0 \\ 1 \end{pmatrix} = (1 \times 0 + 0 \times 1) = \mathbf{0} = (0, 1) \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \langle \mathbf{1} | \mathbf{0} \rangle$$

Moreover

$$\langle \mathbf{0} | \mathbf{0} \rangle = (1, \ 0) \begin{pmatrix} 1 \\ 0 \end{pmatrix} = (1 \times 1 + 0 \times 0) = \mathbf{1} = (0, \ 1) \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \langle \mathbf{1} | \mathbf{1} \rangle$$





Exercises inner product

- Let $|\psi\rangle=\frac{1}{\sqrt{2}}|0\rangle+\frac{1}{\sqrt{2}}|1\rangle$ and $|\phi\rangle=\frac{\sqrt{3}}{2}|0\rangle+\frac{1}{2}|1\rangle$. Compute $\langle\psi|\phi\rangle$ and $\langle\phi|\psi\rangle$
- 2 Let $|\psi\rangle=\frac{i}{\sqrt{2}}|0\rangle+\frac{i}{\sqrt{2}}|1\rangle$ and $|\phi\rangle=\frac{3}{4}|0\rangle+\frac{\sqrt{7}i}{4}|1\rangle$. Compute $\langle\psi|\phi\rangle$ and $\langle\phi|\psi\rangle$.





$$\begin{array}{lll} \textbf{1.} & |\psi\rangle = \frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle \text{ and } |\phi\rangle = \frac{\sqrt{3}}{2}|0\rangle + \frac{1}{2}|1\rangle \\ & \langle\psi|\phi\rangle & = & (\frac{1}{\sqrt{2}}\langle 0| + \frac{1}{\sqrt{2}}\langle 1|)(\frac{\sqrt{3}}{2}|0\rangle + \frac{1}{2}|1\rangle) \\ & = & \frac{1}{\sqrt{2}}\times\frac{\sqrt{3}}{2}\langle 0|0\rangle + \frac{1}{\sqrt{2}}\times\frac{1}{2}\langle 0|1\rangle + \frac{1}{\sqrt{2}}\times\frac{\sqrt{3}}{2}\langle 1|0\rangle + \frac{1}{\sqrt{2}}\times\frac{1}{2}\langle 1|1\rangle \\ & = & \frac{1}{\sqrt{2}}\times\frac{\sqrt{3}}{2}\times 1 + \frac{1}{\sqrt{2}}\times\frac{1}{2}\times 0 + \frac{1}{\sqrt{2}}\times\frac{\sqrt{3}}{2}\times 0 + \frac{1}{\sqrt{2}}\times\frac{1}{2}\times 1 \\ & = & \frac{\sqrt{3}+1}{2\sqrt{2}} \end{array}$$





$$\begin{array}{lll} \textbf{1.} & |\psi\rangle = \frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle \text{ and } |\phi\rangle = \frac{\sqrt{3}}{2}|0\rangle + \frac{1}{2}|1\rangle \\ & \langle \phi|\psi\rangle & = & (\frac{\sqrt{3}}{2}\langle 0| + \frac{1}{2}\langle 1|)(\frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle) \\ & = & \frac{\sqrt{3}}{2} \times \frac{1}{\sqrt{2}}\langle 0|0\rangle + \frac{\sqrt{3}}{2} \times \frac{1}{\sqrt{2}}\langle 0|1\rangle + \frac{1}{2} \times \frac{1}{\sqrt{2}}\langle 1|0\rangle + \frac{1}{2} \times \frac{1}{\sqrt{2}}\langle 1|1\rangle \\ & = & \frac{\sqrt{3}}{2} \times \frac{1}{\sqrt{2}} \times 1 + \frac{\sqrt{3}}{2} \times \frac{1}{\sqrt{2}} \times 0 + \frac{1}{2} \times \frac{1}{\sqrt{2}} \times 0 + \frac{1}{2} \times \frac{1}{\sqrt{2}} \times 1 \\ & = & \frac{\sqrt{3}+1}{2\sqrt{2}} \end{array}$$





$$\begin{aligned} \mathbf{2.} \ |\psi\rangle &= \frac{i}{\sqrt{2}} |0\rangle + \frac{i}{\sqrt{2}} |1\rangle \text{ and } |\phi\rangle = \frac{3}{4} |0\rangle + \frac{\sqrt{7}i}{4} |1\rangle. \\ \langle\psi|\phi\rangle &= \left(\frac{-i}{\sqrt{2}} \langle 0| + \frac{-i}{\sqrt{2}} \langle 1|\right) (\frac{3}{4} |0\rangle + \frac{\sqrt{7}i}{4} |1\rangle) \\ &= \frac{-i}{\sqrt{2}} \times \frac{3}{4} \langle 0|0\rangle + \frac{-i}{\sqrt{2}} \times \frac{\sqrt{7}i}{4} \langle 0|1\rangle + \frac{-i}{\sqrt{2}} \times \frac{3}{4} \langle 1|0\rangle + \frac{-i}{\sqrt{2}} \times \frac{\sqrt{7}i}{4} \langle 1|1\rangle \\ &= \frac{-i}{\sqrt{2}} \times \frac{3}{4} \times 1 + \frac{-i}{\sqrt{2}} \times \frac{\sqrt{7}i}{4} \times 0 + \frac{-i}{\sqrt{2}} \times \frac{3}{4} \times 0 + \frac{-i}{\sqrt{2}} \times \frac{\sqrt{7}i}{4} \times 1 \\ &= \frac{\sqrt{7}}{4\sqrt{2}} - \frac{3}{4\sqrt{2}} \mathbf{i} \end{aligned}$$





2.
$$|\psi\rangle=\frac{i}{\sqrt{2}}|0\rangle+\frac{i}{\sqrt{2}}|1\rangle$$
 and $|\phi\rangle=\frac{3}{4}|0\rangle+\frac{\sqrt{7}i}{4}|1\rangle$.

$$\begin{split} \langle \phi | \psi \rangle &= \left(\frac{3}{4} \langle 0 | + \frac{-\sqrt{7}i}{4} \langle 1 | \right) \left(\frac{i}{\sqrt{2}} | 0 \rangle + \frac{i}{\sqrt{2}} | 1 \rangle \right) \\ &= \frac{3}{4} \times \frac{i}{\sqrt{2}} \langle 0 | 0 \rangle + \frac{3}{4} \times \frac{i}{\sqrt{2}} \langle 0 | 1 \rangle + \frac{-\sqrt{7}i}{4} \times \frac{i}{\sqrt{2}} \langle 1 | 0 \rangle + \frac{-\sqrt{7}i}{4} \times \frac{i}{\sqrt{2}} \langle 1 | 1 \rangle \\ &= \frac{3}{4} \times \frac{i}{\sqrt{2}} \times 1 + \frac{3}{4} \times \frac{i}{\sqrt{2}} \times 0 + \frac{-\sqrt{7}i}{4} \times \frac{i}{\sqrt{2}} \times 0 + \frac{-\sqrt{7}i}{4} \times \frac{i}{\sqrt{2}} \times 1 \\ &= \frac{\sqrt{7}}{4\sqrt{2}} + \frac{3}{4\sqrt{2}} \mathbf{i} \end{split}$$





Linear operator

Def. Linear operator. A linear operator between vector spaces \mathbb{V} and \mathbb{W} is defined as any function $\hat{A}:\mathbb{V}\to\mathbb{W}$ which is linear in its inputs,

$$\hat{A}\left(\sum_{i}\alpha_{i}|\psi_{i}\rangle\right) = \sum_{i}\alpha_{i}\hat{A}|\psi_{i}\rangle$$





Adjoint/Hermitian Conjugate Operator

Let $\hat{A}: \mathcal{H} \to \mathcal{H}$ be a linear operator that induces the map $|\psi\rangle \to |\psi'\rangle$.

The operator \hat{A}^{\dagger} , known as \hat{A} dagger, the adjoint of \hat{A} or the Hermitian Conjugate of \hat{A} , induces the map $\langle \psi | \to \langle \psi' |$ on the corresponding bras.

In other words,

$$\hat{A}|\psi\rangle = |\psi'\rangle$$
$$\langle\psi|\hat{A}^{\dagger} = \langle\psi'|$$

In matrix notation, \hat{A}^{\dagger} is $(A^t)^*$ where t denotes transposition and * denotes complex conjugation.





Hermitian and Unitary Operators

Hermitian operator. Let $\mathcal H$ be a Hilbert space and $\hat A:\mathcal H\to\mathcal H$ a linear operator. If $\hat A=\hat A^\dagger$ then $\hat A$ is a **Hermitian operator**.

Unitary operator. Let $\mathcal H$ be a Hilbert space and $\hat U:\mathcal H\to\mathcal H$ a linear operator. $\hat U$ is a **Unitary operator** if $\hat U\hat U^\dagger=\hat I$, where $\hat I$ is the identity operator.





Outer product

We can also use the Dirac notation to compute vectors. Let $|\psi\rangle, |a\rangle \in \mathcal{H}_1$ and $|\phi\rangle \in \mathcal{H}_2$ then the *outer product* is the linear operator from \mathcal{H}_1 to \mathcal{H}_2 defined by

$$(|\phi\rangle\langle\psi|)|a\rangle \equiv (\langle\psi|a\rangle)|\phi\rangle$$

As it may be expected, the summation of outer products is also a linear operator.





Example - Outer product

For example, let us define the Hadamard operator

$$\hat{H} = \frac{1}{\sqrt{2}}(|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| - |1\rangle\langle 1|)$$

The action of \hat{H} on ket $|0\rangle$ is given by

$$\begin{split} \hat{H}|\mathbf{0}\rangle &= \left(\frac{1}{\sqrt{2}}|0\rangle\langle 0| + \frac{1}{\sqrt{2}}|0\rangle\langle 1| + \frac{1}{\sqrt{2}}|1\rangle\langle 0| - \frac{1}{\sqrt{2}}|1\rangle\langle 1|\right)|\mathbf{0}\rangle \\ &= \frac{\langle 0|\mathbf{0}\rangle}{\sqrt{2}}|0\rangle + \frac{\langle 1|\mathbf{0}\rangle}{\sqrt{2}}|0\rangle + \frac{\langle 0|\mathbf{0}\rangle}{\sqrt{2}}|1\rangle - \frac{\langle 1|\mathbf{0}\rangle}{\sqrt{2}}|1\rangle \\ &= \frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle \end{split}$$





Exercise - Outer product

Let
$$\hat{\sigma}_y=i|0\rangle\langle 1|-i|1\rangle\langle 0|$$
 and $|\psi\rangle=\frac{\sqrt{3}}{2}|0\rangle+\frac{i}{2}|0\rangle$. Compute $\hat{\sigma}_y|\psi\rangle$.





Answer to exercise - Outer product

Let $\hat{\sigma}_y=i|0\rangle\langle 1|-i|1\rangle\langle 0|$ and $|\psi\rangle=\frac{\sqrt{3}}{2}|0\rangle+\frac{i}{2}|1\rangle$. Compute $\hat{\sigma}_y|\psi\rangle$.

$$\hat{\sigma}_{y}|\psi\rangle = (i|0\rangle\langle 1| - i|1\rangle\langle 0|)(\frac{\sqrt{3}}{2}|0\rangle + \frac{i}{2}|1\rangle)
= \frac{\sqrt{3}i\langle 1|0\rangle}{2}|0\rangle + \frac{i^{2}\langle 1|1\rangle}{2}|0\rangle - \frac{\sqrt{3}i\langle 0|0\rangle}{2}|1\rangle - \frac{i^{2}\langle 0|1\rangle}{2}|1\rangle
= -\frac{1}{2}|0\rangle - \frac{\sqrt{3}i}{2}|1\rangle$$





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In quantum computing, information is stored, manipulated and measured in the form of qubits.

A qubit may be mathematically represented as a unit vector in a two-dimensional Hilbert space $|\psi\rangle\in\mathcal{H}^2.$

A qubit $|\psi\rangle$ may be written in general form as

$$|\psi\rangle = \alpha|p\rangle + \beta|q\rangle$$

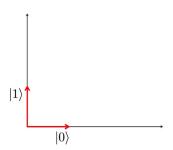
where $\alpha, \beta \in \mathbb{C}$, $|\alpha|^2 + |\beta|^2 = 1$ and $\{|p\rangle, |q\rangle\}$ is an arbitrary basis spanning \mathcal{H}^2 .





A most important consequence of the vectorial nature of a qubit is the possibility of writing it as a linear combination of elements of any basis.

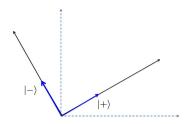
For example, we may choose $|p\rangle=|0\rangle$ and $|q\rangle=|1\rangle$ to write a qubit. The basis $\{|0\rangle,|1\rangle\}$ is known as **the computational basis**.







We may also choose the **the diagonal basis** $\{|+\rangle, |-\rangle\}$ to write a qubit.







Choosing a concrete vector basis and values for corresponding complex coefficients is known as **preparing a qubit**. For instance,

$$|\psi\rangle = \frac{\sqrt{3}}{2}|0\rangle + \frac{i}{2}|1\rangle$$

In computer science parlance, preparing a qubit (or a set of qubits) is equivalent to variable initialization in a computer program.





Postulate 2. Evolution by Unitary Operator

The evolution of a closed quantum system with state vector $|\Psi\rangle$ is described by a Unitary operator. The state of a system at time t_2 according to its state at time t_1 is given by

$$|\Psi(t_2)\rangle = \hat{U}|\Psi(t_1)\rangle.$$

Postulate 2 only describes the mathematical properties that an evolution operator must have. The specific evolution operator required to describe the behaviour of a particular quantum system depends on the system itself.





Postulate 2. Evolution by Schrödinger equation

Quantum evolution can also be written in terms of differential equations.

The time evolution of a closed quantum system can be described by the Schrödinger equation:

$$i\hbar \frac{\partial |\psi(t)\rangle}{\partial t} = \mathbf{\hat{H}}|\psi(t)\rangle$$

where $\hbar = \frac{h}{\pi}$, h is Plank Constant, and $\hat{\mathbf{H}}$ is a Hermitian operator known as the *Hamiltonian* of the sytem.





Examples of Unitary Operators (1/2)

Pauli operators

$$\hat{\sigma}_x = |0\rangle\langle 1| + |1\rangle\langle 0|; \hat{\sigma}_y = i|0\rangle\langle 1| - i|1\rangle\langle 0|; \hat{\sigma}_z = |0\rangle\langle 0| - |1\rangle\langle 1|$$

If
$$|0\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix}$$
, $|1\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$, $\langle 0| = (1, 0)$, and $\langle 1| = (0, 1)$

Then we produce the following matrix representations of Pauli operators:

$$\sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$
 ; $\sigma_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}$; $\sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$





Examples of Unitary Operators (2/2)

The **Hadamard operator**

$$\hat{H} = \frac{1}{\sqrt{2}}(|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| - |1\rangle\langle 1|)$$

Again, if
$$|0\rangle=\begin{pmatrix}1\\0\end{pmatrix}$$
, $|1\rangle=\begin{pmatrix}0\\1\end{pmatrix}$, $\langle 0|=(1,\ 0)$, and $\langle 1|=(0,\ 1)$

$$H = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1\\ 1 & -1 \end{pmatrix}$$





(Already known) Example of Evolution by Unitary Operator

Let $\hat{\sigma}_y=i|0\rangle\langle 1|-i|1\rangle\langle 0|$ and $|\psi\rangle=\frac{\sqrt{3}}{2}|0\rangle+\frac{i}{2}|1\rangle.$ Compute $\hat{\sigma}_y|\psi\rangle.$

$$\begin{array}{lll} \hat{\sigma}_{y}|\psi\rangle & = & (i|0\rangle\langle1|-i|1\rangle\langle0|)(\frac{\sqrt{3}}{2}|0\rangle + \frac{i}{2}|1\rangle) \\ \\ & = & \frac{\sqrt{3}i\langle1|0\rangle}{2}|0\rangle + \frac{i^{2}\langle1|1\rangle}{2}|0\rangle - \frac{\sqrt{3}i\langle0|0\rangle}{2}|1\rangle - \frac{i^{2}\langle0|1\rangle}{2}|1\rangle \\ \\ & = & -\frac{1}{2}|0\rangle - \frac{\sqrt{3}i}{2}|1\rangle \end{array}$$





In quantum mechanics, measurement is a non-trivial and highly counter-intuitive process because:

a) Measurement outcomes are inherently probabilistic. Regardless of the carefulness in the preparation of a measurement procedure, the possible outcomes of such measurement are distributed according to a certain probability distribution.





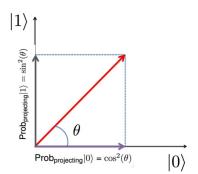
b) Pre- and post-measurement quantum states are different.

Once a measurement has been performed, a quantum system is unavoidably altered due to the interaction with the measurement apparatus. Consequently, for an arbitrary quantum system, pre-measurement and post-measurement quantum states are different in general.





c) Measuring \Leftrightarrow Projecting. Measuring a quantum system is equivalent to projecting (as in analytic geometry) a vector onto a vector basis.









The **KEY** points to remember are:

- Measuring a quantum system is a probabilistic process that changes the state of the quantum system.
 - Extracting information from a quantum system (e.g. a photon) makes irreversible changes in the actual information contained in that quantum system.
 - Measuring a quantum system is equivalent to projecting a vector onto a vector basis.





Example. Let $|\psi\rangle$ be a qubit representing a two-dimensional quantum-mechanical system (e.g., photon polarization). Let us write $|\psi\rangle$ as

$$|\psi\rangle = \alpha|0\rangle + \beta|1\rangle$$

where $||\alpha||^2+||\beta||^2=1$ and $\{|0\rangle,|1\rangle\}$ is the computational basis spanning \mathcal{H}^2 .





According to Postulate 3, the probabilities of obtaining outcome a_0 or outcome a_1 are given by

$$p(a_0) = ||\alpha||^2$$
 and $p(a_1) = ||\beta||^2$

Corresponding post-measurement quantum states are as follows:

If outcome =
$$a_0$$
 then $|\psi\rangle_{pm} = |0\rangle$
If outcome = a_1 then $|\psi\rangle_{pm} = |1\rangle$.





We now focus on the mathematical description of a composite quantum system, i.e. a system made up of several different physical systems.

The state space of a composite quantum system is the tensor product of the component system state spaces.

If we have n quantum systems expressed as *state vectors*, labeled $|\psi\rangle_1, |\psi\rangle_2, \ldots, |\psi\rangle_n$ then the joint state of the total system is given by $|\psi\rangle_T = |\psi\rangle_1 \otimes |\psi\rangle_2 \otimes \ldots \otimes |\psi\rangle_n$.

Please note this crucial property: $\otimes_{i=1}^n \mathcal{H}_i^2 = \mathcal{H}^{2^n}$.





Let $\hat{A}:\mathcal{H}_1\to\mathcal{H}_1$ and $\hat{B}:\mathcal{H}_2\to\mathcal{H}_2$ be linear operators. Then, the following equalities hold $\forall\ |a\rangle_1,|a\rangle_2\in\mathcal{H}_1,\ |b\rangle_1,|b\rangle_2\in\mathcal{H}_2,$ $\alpha\in\mathbb{C}$:

1)
$$\alpha(|a\rangle_1\otimes|b\rangle_1)=(\alpha|a\rangle_1)\otimes|b\rangle_1=|a\rangle_1\otimes(\alpha|b\rangle_1)$$

2)
$$(|a\rangle_1 + |a\rangle_2) \otimes |b\rangle_1) = |a\rangle_1 \otimes |b\rangle_1 + |a\rangle_2 \otimes |b\rangle_1$$

3)
$$|a\rangle_1 \otimes (|b\rangle_1 + |b\rangle_2) = |a\rangle_1 \otimes |b\rangle_1 + |a\rangle_1 \otimes |b\rangle_2$$

4)
$$\hat{A} \otimes \hat{B}(|a\rangle_1 \otimes |b\rangle_1) = \hat{A}|a\rangle_1 \otimes \hat{B}|b\rangle_1$$

5) Let
$$|a\rangle_i \in \mathcal{H}_1$$
, $|b\rangle_i \in \mathcal{H}_2$ and $\alpha_i \in \mathbb{C} \Rightarrow \hat{A} \otimes \hat{B}(\sum_i \alpha_i |a\rangle_i \otimes |b\rangle_i) = \sum_i \alpha_i \hat{A} |a\rangle_i \otimes \hat{B} |b\rangle_i$





- We may write $|a\rangle \otimes |b\rangle$ as $|ab\rangle$ or $|a,b\rangle$.
- Moreover, the tensor product of $|a\rangle$ with itself n times may be written as $|a\rangle\otimes|a\rangle\otimes\ldots\otimes|a\rangle=|a\rangle^{\otimes n}$.





The Kronecker product is a matrix representation of the tensor product and it is defined as follows.

Let $A=(a_{ij}), B=(b_{ij})$ be two matrices of order $m\times n$ and $p\times q$ respectively. Then $A\otimes B$ is given by

$$A \otimes B = \begin{pmatrix} A_{11}B & A_{12}B & \dots & A_{1n}B \\ A_{21}B & A_{22}B & \dots & A_{2n}B \\ \vdots & \vdots & \vdots & \vdots \\ A_{m1}B & A_{m2}B & \dots & A_{mn}B \end{pmatrix}.$$

 $A\otimes B$ is of order $mp\times nq$.





Example 1. Let $\hat{H}^{\otimes 2}$ be the tensor product of the Hadamard operator with itself and let $|\psi\rangle=|0\rangle\otimes|0\rangle=|00\rangle$. Then

$$\begin{array}{ll} \hat{H}^{\otimes 2} \mid \pmb{\psi} \rangle & = & \frac{1}{2} (\mid \! 00 \rangle \langle 00 \mid + \mid \! 01 \rangle \langle 00 \mid + \mid \! 10 \rangle \langle 00 \mid + \mid \! 11 \rangle \langle 00 \mid + \mid \! 00 \rangle \langle 01 \mid \\ & - \mid \! 01 \rangle \langle 01 \mid + \mid \! 10 \rangle \langle 01 \mid - \mid \! 11 \rangle \langle 01 \mid + \mid \! 00 \rangle \langle 10 \mid + \mid \! 01 \rangle \langle 10 \mid \\ & - \mid \! 10 \rangle \langle 10 \mid - \mid \! 11 \rangle \langle 10 \mid + \mid \! 00 \rangle \langle 11 \mid - \mid \! 01 \rangle \langle 11 \mid - \mid \! 10 \rangle \langle 11 \mid \\ & + \mid \! 11 \rangle \langle 11 \mid) \mid \! 00 \rangle \\ & = & \frac{1}{2} (\langle 00 \mid \! 00 \rangle \mid \! 00 \rangle + \langle 00 \mid \! 00 \rangle \mid \! 01 \rangle + \langle 00 \mid \! 00 \rangle \mid \! 10 \rangle + \langle 00 \mid \! 00 \rangle \mid \! 11 \rangle \\ & = & \frac{1}{2} (\mid \! 00 \rangle + \mid \! 01 \rangle + \mid \! 10 \rangle + \mid \! 11 \rangle) \end{array}$$

That is, the notion of inner product on orthonormal bases easily extends to \mathcal{H}^{2^n} .





Example 2. Introducing \hat{C}_{not} .

Let us introduce \hat{C}_{not} , a two-qubit gate that behaves as follows:

$$\hat{C}_{\mathsf{not}} = |00\rangle\langle00| + |01\rangle\langle01| + |10\rangle\langle11| + |11\rangle\langle10|$$

Then

$$\hat{C}_{\mathsf{not}}|00\rangle = |00\rangle$$

$$\hat{C}_{\mathsf{not}}|01\rangle = |01\rangle$$

$$\hat{C}_{\mathsf{not}}|10\rangle = |11\rangle$$

$$\hat{C}_{\mathsf{not}}|11\rangle = |10\rangle$$

The operator \hat{C}_{not} acts as a *Controlled Not*: \hat{C}_{not} flips the second qubit iff the first qubit is $|1\rangle$, otherwise \hat{C}_{not} behaves like a buffer.





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Introduction to Quantum Circuits

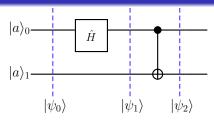
We now introduce a quantum circuit to compute Bell states







Bell State Circuit (1/4)

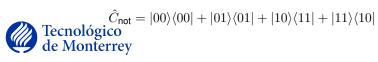


Let

$$|a\rangle_0=|0\rangle$$
 and $|a\rangle_1=|0\rangle$

Also, remember that

$$\hat{H} = \frac{1}{\sqrt{2}}(|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| - |1\rangle\langle 1|)$$





Bell State Circuit (2/4)

So,

$$\begin{split} |\psi\rangle_0 &= |0\rangle\otimes|0\rangle = |00\rangle \\ |\psi\rangle_1 &= (\hat{H}\otimes\hat{\mathbb{I}})(|0\rangle\otimes|0\rangle) = \hat{H}|0\rangle\otimes\hat{\mathbb{I}}|0\rangle = \frac{1}{\sqrt{2}}(|0\rangle+|1\rangle)\otimes|0\rangle \\ &= \frac{1}{\sqrt{2}}|0\rangle|0\rangle + \frac{1}{\sqrt{2}}|1\rangle|0\rangle \\ |\psi\rangle_2 &= \hat{C}_{\mathrm{not}}(\frac{1}{\sqrt{2}}|0\rangle|0\rangle + \frac{1}{\sqrt{2}}|1\rangle|0\rangle) \\ &= \frac{1}{\sqrt{2}}|0\rangle|0\rangle + \frac{1}{\sqrt{2}}|1\rangle|1\rangle = \frac{1}{\sqrt{2}}|00\rangle + \frac{1}{\sqrt{2}}|11\rangle \end{split}$$

So,

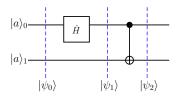


$$|\psi\rangle_2 = \frac{|00\rangle + |11\rangle}{\sqrt{2}}$$



Bell State Circuit (3/4)

Exercise



Compute $|\psi\rangle_2$ for

$$\begin{array}{l} |a\rangle_0=|0\rangle \text{ and } |a\rangle_1=|1\rangle \\ |a\rangle_0=|1\rangle \text{ and } |a\rangle_1=|0\rangle \\ |a\rangle_0=|1\rangle \text{ and } |a\rangle_1=|1\rangle \end{array}$$





Bell State Circuit (4/4)

Answers

$$\begin{split} \hat{C}_{\mathsf{not}}((\hat{H}\otimes\hat{\mathbb{I}})|00\rangle) &= \frac{|00\rangle + |11\rangle}{\sqrt{2}} \\ \hat{C}_{\mathsf{not}}((\hat{H}\otimes\hat{\mathbb{I}})|01\rangle) &= \frac{|01\rangle + |10\rangle}{\sqrt{2}} \\ \hat{C}_{\mathsf{not}}((\hat{H}\otimes\hat{\mathbb{I}})|10\rangle) &= \frac{|00\rangle - |11\rangle}{\sqrt{2}} \\ \hat{C}_{\mathsf{not}}((\hat{H}\otimes\hat{\mathbb{I}})|11\rangle) &= \frac{|01\rangle - |10\rangle}{\sqrt{2}} \end{split}$$

These states are known as the Bell states

$$\begin{split} |\Phi^{+}\rangle &= \frac{|00\rangle + |11\rangle}{\sqrt{2}} \\ |\Phi^{-}\rangle &= \frac{|00\rangle - |11\rangle}{\sqrt{2}} \\ |\Psi^{+}\rangle &= \frac{|01\rangle + |10\rangle}{\sqrt{2}} \\ |\Psi^{-}\rangle &= \frac{|01\rangle - |10\rangle}{\sqrt{2}} \end{split}$$





Quantum Entanglement (1/2)

Bell states are examples of entangled states. Bell states are key features of a quantum information transmission protocol known as quantum teleportation.

Quantum entanglement is a unique type of correlation shared between components of a quantum system.

Quantum entanglement and the principle of superposition are two of the main features behind the power of quantum computation and quantum information theory.





Quantum Entanglement (2/2)

Entangled quantum systems are sometimes best used collectively, that is, sometimes an optimal use of entangled quantum systems for information storage and retrieval includes manipulating and measuring those systems as a whole, rather than on an individual basis.





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General structure of a quantum algorithm

	Classical algorithm	Quantum Algorithm
Information Unit	Bit $x \in \{0, 1\}$	Qubit $ \psi angle\in\mathbb{C}^2$
Algorithm design	{AND, OR, NOT}	$ \Psi_{t_2}\rangle = \hat{U} \Psi_{t_1}\rangle$
	Automata (e.g. Turing Machines)	$i\hbar rac{d \psi\rangle}{dt} = \mathbf{\hat{H}} \psi\rangle$
	C, Python	©
Outcome	Memory readout	Quantum measurement





How to write quantum algorithms

- State of the art on quantum programming languages:
 - Quantum gate-based graphical interface IBM.
 - Python, C, C++ plus APIs D-Wave
 - Full quantum programming languages: yet to come
- Software for digital simulation of quantum algorithms
 - IBM. Q interface and QSKIT (Python)
 - Microsoft. Liquid and Q#
 - D-Wave. SAPI.





How to write quantum algorithms

There are several methods to write a quantum algorithm.

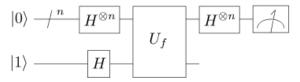






Quantum gates

A most popular method to write a quantum algorithm is to use quantum gates/circuits, which are mathematical representations of physical transformations on quantum systems.



Deutsch-Jozsa quantum algorithm. Source: Wikipedia ®

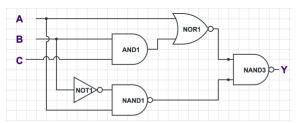
This programming method is based on $|\Psi_{t_2}\rangle = \hat{U}|\Psi_{t_1}\rangle$.





Quantum gates

The rationale behind the quantum gate method is very similar to the one followed in the construction of digital circuits:



A digital circuit employing NAND, NOR, NOT and AND gates. Source: Wikipedia ®





Quantum gates

- The mathematical description of digital circuitry is determined and constrained by the definitions and results of mathematical logic.
- As for quantum circuits, the mathematical description of a quantum gate is determined by the physics of quantum evolution (that is, the behaviour of quantum systems in time).





Schrödinger equation

We may also use differential equations to write quantum algorithms:

$$i\hbar \frac{d|\psi\rangle}{dt} = \hat{\mathbf{H}}|\psi\rangle$$

This is the Schrödinger equation and it is basic equation in the models of quantum annealing and adiabatic quantum computation.





Example - Grover's Algorithm

A quantum algorithm developed by Lov Grover, its purpose is to perform a search of an item from an **unsorted and unstructured** list of N records.

- A classical computer brute-force *sequential* algorithm would require O(N) steps to perform the search.
- In contrast, Grover's algorithm requires only $O(\sqrt{N})$ steps.
- Key concepts: quantum superposition, quantum parallelism and quantum oracle-based computation.





Example - Shor's Algorithm

A quantum algorithm developed by Peter Shor, its purpose is to efficiently compute the prime factors of an N-digit integer ($N=2^n$).

- A classical computer brute-force algorithm would require $O(2^{n/2})$ steps to perform the search.
- Quantum solution: Shor's algorithm, $O(n^3)$ steps.
- Key concept: Shor's algorithm provides exponential acceleration wrt known classical algorithms.



