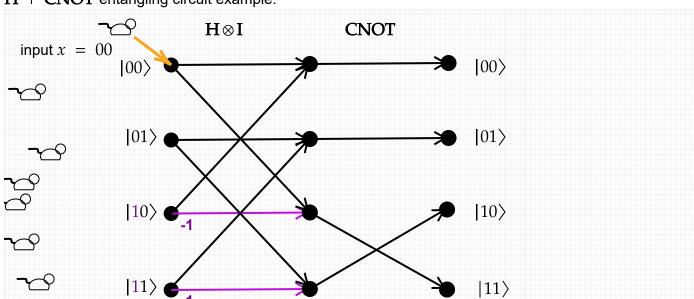
CSE491/596 Lecture Fri. 12/10/21: Visualizing Small-Scale Quantum Applications

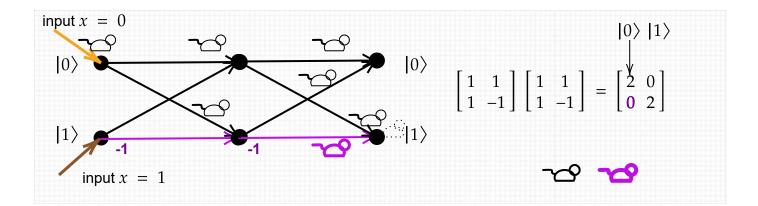
For the textbook, I devised a "maze" visualization of quantum **amplification** and **interference** for applications with up to 3 or so qubits. I found that the brilliant Dorit Aharonov had used the same idea. Most basic gate matrices have the following "balance" property: all nonzero entries have the same magnitude, so normalizing factors like $\frac{1}{\sqrt{2}}$ can be set aside. Then the Hadamard, **CNOT**, Toffoli, and Pauli gates (ignoring the global *i* factor in **Y**) give just entries +1 or -1, which become the only possible values of any *Feynman path*, which means a multiplicative term in the ultimate matrix product. That makes it easier to sum the results of paths. The index values m, l, k, j, i, \ldots become "locations" in the wavefront as it flows for time *s*. Since the paths are discrete we can picture "lab mice" running through the maze. The humorous treatment in section 7.5 illustrates:

- 1. superposition
- 2. interference
- 3. amplification
- 4. measurement
- 5. the system's quantum state after a measurement.



One nice thing is that you can read the mazes left-to-right, same as the circuits. Here is the H + CNOT entangling circuit example:

No interference or amplification is involved here---the point is that if you enter at $|00\rangle$, then $|00\rangle$ and $|11\rangle$ are the only places you can come out---and they have equal weight. To see interference, you can string the "maze gadgets" for two Hadamard gates together:



In linear-algebra terms, all that happened at lower right was $1 \cdot 1 + -1 \cdot 1$ giving 0. But the wave interference being described that way is a real physical phenomenon. Even more, according to Deutsch the two serial Hadamard gates branch into 4 universes, each with its own "Phil the mouse" (which can be a photon after going through a beam-splitter). One of those universes has "Anti-Phil", who attacks a "Phil" that tries to occupy the same location (coming from a different universe) and they fight to mutual annihilation.

Examples in Chapter 8, "Deutsch's Algorithm"

David Deutsch, drawing on two papers by Feynman and other sources, introduced quantum computing while he and I were graduate students at Oxford in the mid-1980s. At first, he claimed quantum computers could solve the Halting Problem in finite time. Fellows of Oxford's Mathematical Institute refuted the claim. But it was not crazy: a year ago it was proved that a binary quantum system of "interactive provers" *can* (kind-of-)solve the Halting Problem in finite time. (My review of the paper is at https://rjlipton.wordpress.com/2020/01/15/halting-is-poly-time-quantum-provable/) Per my memory of observing some meetings about it, the gap in Deutsch's argument had to do with properties of probability measures based on infinite binary sequences.

So Deutsch fell back on something less ambitious: demonstrating that there was a "very finite" task that quantum computers can do and classical ones cannot. (Well, unless the playing field is leveled for them...but before we argue about it, let's see the task.) The task is a **learning problem**, a kind of interaction we haven't covered until this last day. Instead of "input *x*, compute y = f(x)", a learning problem is to determine facts about an initially-unknown entity *f* that you can **query**.

1. Oracle Turing machines give a classic way to define this kind of problem. For oracle functions f or languages A drawn from a limited class---such as subclasses of the regular languages--can we design an OTM M that on input 0^n (for large enough n) can distinguish what A is in time (say) polynomial in n? The computation $M^A(0^n)$ can learn about A by making queries y on selected strings y and observing the answers A(y).

- 2. One can also define **oracle circuits** that have special **oracle gates** with some number *m* of input wires and enough output wires to give the answer f(y) on any $y \in \{0, 1\}^m$.
- 3. An ordinary electrical test kit behaves that way. It is a circuit with a place(s) for you to insert one or more (possibly-defective) electrical components *A*. The test results should diagnose electrical facts about *A*.
- 4. Quantum circuits for all of the Deutsch, Deutsch-Jozsa, Simon, Shor, and Grover algorithms work this way. They involve an **oracle function** $f: \{0,1\}^n \rightarrow \{0,1\}^r$ given in **reversible form** as the function $F: \{0,1\}^{n+r} \rightarrow \{0,1\}^{n+r}$ defined by:

$$F(x,z) = (x, f(x) \oplus z).$$

Usually *z* is 0^r and the comma is just concatenation (i.e., tensor product) so the output is just xf(x). In the simplest case n = r = 1, *F* is a two-(qu)bit function. Some examples:

- If f is the identity function, f(x) = x, then $F(x, z) = (x, x \oplus z) = \text{CNOT}(x, z)$.
- If $f(x) = \neg x$, then $F(x, z) = (x, x \leftrightarrow z)$: F(00) = 01, F(01) = 00, F(10) = 10, F(11) = 11.
- If f is always false, i.e., f(x) = 0, then F is the identity function.
- If f(x) = 1, then $F(x, z) = (x, \neg z)$, so F(00) = 01, F(01) = 00, F(10) = 11, F(11) = 10.

These are all deterministic as functions of two-qubit basis states, so they permute the quantum coordinates 1 = 00, 2 = 01, 3 = 10, and 4 = 11. Recall that **CNOT** gives the permutation that swaps the coordinates 3 and 4, that is, **CNOT** = (3 4) in swap notation. In full, we have:

$$F_{id} = (34), F_{\neg} = (12), F_{0} = (), F_{1} = (12)(34).$$

The functions f(x) = 0 and f(x) = 1 are *constant*. The identity and \neg functions have one true and one false value each, so they *balance* values of 0 and 1. The question posed by Deutsch is:

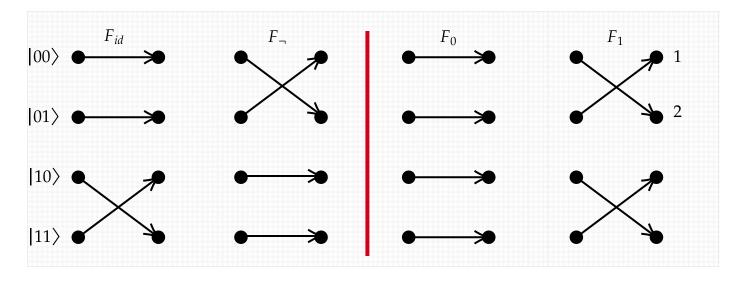
How many queries are needed to tell whether f is constant from whether f is balanced?

If we just think of f, suppose we try the query y = 0 and ask for f(y). If we get the answer " f(0) = 0" then it f could be constant-false, but f could also be the balanced identity function. The answer f(0) = 1 would leave both constant-true and negation as possibilities. Likewise if we try y = 1. The first point is that this impossibility of hitting things with one query carries forward to the way we have to modify the problem for quantum:

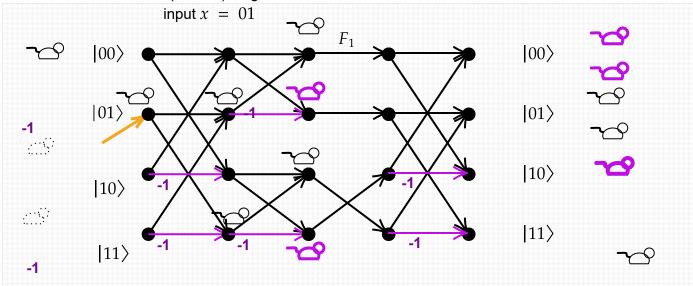
How many queries are needed to tell ($F_{id} \text{ or } F_{\neg}$) apart from ($F_0 \text{ or } F_1$)?

It seems like we have more of a chance because now we can query two things: 00, 01, 10, or 11. Or in the permutation view, we can query y = 1, 2, 3, or 4. The problem is that the range of answers we can get is too limited for this to help. F(1) and F(2) can only be 1 ro 2; F(3) and F(4) can only be 3 or

4. So suppose you query y = 3 and get the answer 4. Then *F* could be F_{id} or *F* could be F_1 . The basic problem for a classical algorithm is that every quadrant of the following diagram has both a straight and a cross:



A quantum circuit, however, can make one query to an oracle gate for any of these four functions, and can distinguish a member of the first pair from a member of the second pair by the answer to one qubit after a measurement. The input is not $|00\rangle$ but instead $|01\rangle$; that is, the ancilla is initialized to 1, not to 0. Here is the wavefront ("maze") diagram of how it works:



There is, IMHO, an "unfair" aspect of the comparison. The classical algorithm is being allowed to evaluate the oracle only at basis vectors. The quantum algorithm gets to evaluate it at a linear combination---indeed, it's the state

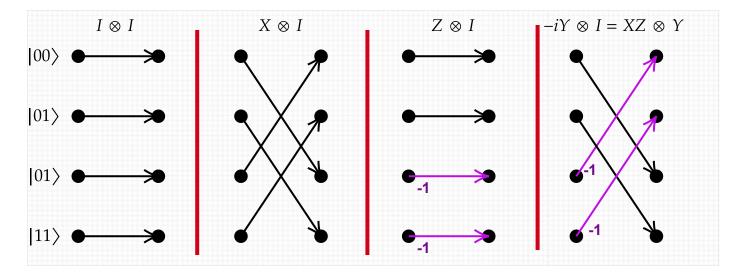
$$|+-\rangle = \frac{1}{2} (|00\rangle - |01\rangle + |10\rangle - |11\rangle)$$

from the Fri. 12/4 lecture. If we do the kind of linear extension of Boolean logic that was covered as the "Binary Linear Equations" presentation option, then we can solve the problem in one shot classically by evaluating at the point (1, -1, 1, -1) and seeing where the – signs end up in the resulting vector. FYI: https://rjlipton.wordpress.com/2011/10/26/quantum-chocolate-boxes/

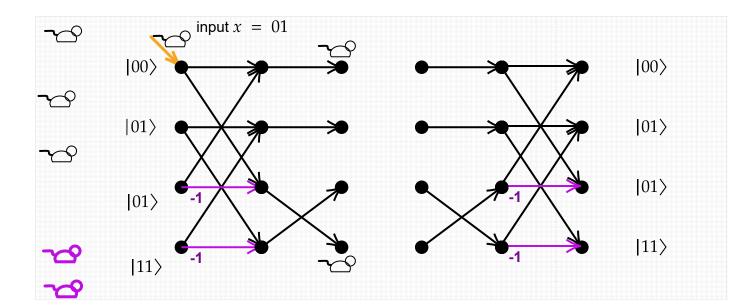
Superdense Coding

It is easy to rig cases F_0 , F_1 , F_2 , F_3 where you can distinguish them exactly by asking one query and measuring both qubits. Just define $F_i(00) = i$, for instance. "Superdense coding" is a case where the rigging has a bit of surprise because it appears to convey 2 bits of information with just 1 qubit of communication. This is impossible by **Holevo's Theorem** that *n* qubits can yield only *n* bits of classical information. (Another instance of this that you can input $\sim \frac{1}{2}n^2$ bits of information by choosing the **CZ** gates for edges of an undirected *n*-vertex graph *G* in a graph-state circuit C_G on *n* qubits, one for each vertex, but you can only get *n* bits of information out by measuring. Hence graph-state encoding is majorly *lossy*.) The rub is that the rigging involves the communicating parties "Alice" and "Bob" already having exchanged 1 bit of information in order to set them up with an entangled qubit pair.

We will regard it as a nice case of the learning problem because it uses the four Pauli matrices. We want to identify one of the following four possibilities **exactly** by the results of **two** qubits.



This time the input is $|00\rangle$. To work it out via wavefronts:



Deutsch-Jozsa Extension

Getting back to Deutsch's Problem, Richard Jozsa added that if you only care about distinguishing *constant* functions $f: \{0,1\}^n \to \{0,1\}$ from *balanced* ones, then you can make the classical algorithms require $2^{n-1} + 1$ queries, while the quantum ones can still do it on one query to a completely separable superposed state. This is a conditional problem, called a **promise problem**, in that it only applies when *f* is in one of those two cases. If *f* is neither balanced nor constant, then "all bets are off"---any answer is fine, even $| - (v) | - \rangle$.

The maze diagrams would get exponentially big, but we can track the computations via linear algebra. It is like Deutsch's setup except with $\mathbf{H}^{\otimes n}$ in place of the first \mathbf{H} , input $|0^n1\rangle$ in place of $|01\rangle$, and targets (ignoring the $\sqrt{2}$ normalizers):

- constant $\mapsto |0^n\rangle(|0\rangle + |1\rangle)$ (instead of $(|00\rangle + |01\rangle)$, so that 0^n is certainly measured.
- balanced $\mapsto |?\rangle$ (instead of $(|10\rangle + |11\rangle)$, such that 0^n is certainly *not* measured.

The key observation is that for any f, any argument $x \in \{0,1\}^n$, and $b \in \{0,1\}$, the amplitude in the component xb of the final quantum state ϕ is

$$\frac{1}{\sqrt{2^{n+1}}} \sum_{t \in \{0,1\}^n} (-1)^{x \bullet t} (-1)^{f(t) \oplus b}.$$

Here $x \bullet t$ means taking the dot-products $x_i \cdot t_i$ (which is the same as $x_i \wedge t_i$) and adding them up modulo 2 (which is the same as XOR-ing them). Well, when $x = 0^n$ this is always just zero, so the first term is $(-1)^0$ and just drops out, leaving

$$\phi(0^n b) = \frac{1}{\sqrt{2^{n+1}}} (-1)^b \sum_{t \in \{0,1\}^n} (-1)^{f(t)}.$$

Note that the $(-1)^{b}$ term is independent of the sum over *t*, so it comes out of the sum---and this is why we get two equal possibilities in the original Deutsch's algorithm as well. Ths final point is that:

- When *f* is *constant*, these terms are all the same, so they *amplify*---giving $\frac{1}{\sqrt{2}}$ for the constant-false function and $\frac{-1}{\sqrt{2}}$ for constant-true. Both of these amplitudes square to $\frac{1}{2}$ and so together soak up all the output probability, so that 0^n is measured with certainty.
- When f is *balanced*, the big sum has an equal number of +1 and -1 terms, so they all *interfere* and *cancel*. Hence 0^n will certainly not be measured.

Added: A *randomized* classical algorithm can efficiently tell with high probability whether f is constant by querying some random strings. If it ever gets different answers $f(y) \neq f(y')$ then definitely f is not constant. (So, under the condition of the "promised problem," it must be balanced.) If it always gets the same answer, then since any balanced function gives 50-50 probability on random strings, it can quickly figure that f is constant. But it is still the case that a deterministic algorithm needs exponentially many queries and hence exponential time.

Brief Conclusion: Simon's Algorithm to Shor's to the Present

Daniel Simon extended Deutsch-Jozsa to a problem where one can prove that a classical randomized algorithm needs exponential time. I still have reservations about whether the classical setting is fair (FYI, see https://rjlipton.wordpress.com/2011/11/14/more-quantum-chocolate-boxes/), but there is no doubt about what it led to. Peter Shor in 1993 realized that if he substituted QFT_n for Deutsch-Jozsa and Simon's use of $H^{\otimes n}$, then he could make it impact the (group-theoretic) periodicity in the powering function modulo a number M = pq, in a way that allows finding p and q with high probability in roughly-quadratic quantum time. That is, he classified **FACTORING** as belonging to BQP. This plus the **de-randomization** of **PRIMES** from RP \cap co-RP to P in 2002 and **UGAP** (which is the graph-accessibility problem for undirected graphs) from randomized logspace to L in 2004 completes our current top-level knowledge about the landscape of major complexity classes:

