# Lecture 2 Computational Hardness

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# 1 Recap: Perfect Secrecy

Recall from last time:

- Shannon secrecy says that the *a posteriori* distribution over the message, given the ciphertext, is identical to the *a priori* distribution. (See the lecture notes for formal definition.) Informally, "seeing the ciphertext is only as good as seeing nothing at all."
- Perfect secrecy is the property that the distribution of the ciphertext  $c \in \mathcal{C}$  (over the choice of key  $k \in \mathcal{K}$ ) is exactly the same, no matter what message  $m \in \mathcal{M}$  was encrypted.
- Shannon secrecy and perfect secrecy are equivalent.
- The one-time pad scheme enjoys Shannon/perfect secrecy.

# 2 Limits of Shannon/Perfect Secrecy

For reference, here again is the definition of Shannon secrecy:

**Definition 2.1** (Shannon secrecy). A shared-key encryption scheme (Gen, Enc, Dec) with message space  $\mathcal{M}$  and ciphertext space  $\mathcal{C}$  is *Shannon secret with respect to a probability distribution* D over  $\mathcal{M}$  if for all  $\bar{m} \in \mathcal{M}$  and all  $\bar{c} \in \mathcal{C}$ ,

$$\Pr_{m \leftarrow D, k \leftarrow \mathsf{Gen}}[m = \bar{m} \mid \mathsf{Enc}_k(m) = \bar{c}] = \Pr_{m \leftarrow D}[m = \bar{m}]. \tag{2.1}$$

The scheme is *Shannon secret* if it is Shannon secret with respect to every distribution D over  $\mathcal{M}$ .

Now that we know that the one-time pad is Shannon secret, what more is left to do? (Cancel the course?) Well, first notice that the one-time pad scheme is not necessarily very usable: the key length is as large as the message length, and for long messages it may be unrealistic for Alice and Bob to establish (and keep secret!) a huge key before communicating. (This is not to mention the additional distinct keys needed for when Carol, Dan, Edith, and Fred enter the picture...) Also notice that the key can by used only *one time*: encrypting more than one message with the same key is "abusing the model," and causes the security proof to break down. In fact, using the key more than once often makes the scheme *completely insecure*. (The USSR learned this the hard way as a result of the US/UK-led "Venona" project during the Cold War.)

It turns out that these drawbacks are not really a deficiency of the one-time pad itself, but are *inherent* in the strong requirement of perfect secrecy.

**Theorem 2.2** (Shannon's theorem). *If a shared-key encryption scheme (with key space* K *and message space* M) *is Shannon secret, then*  $|K| \ge |M|$ .

*Proof.* We prove the contrapositive: supposing that  $|\mathcal{K}| < |\mathcal{M}|$  for some scheme (Gen, Enc, Dec), we will show that the scheme is not Shannon secret. It suffices to show the existence of a distribution D over  $\mathcal{M}$ , a fixed message  $\bar{m} \in \mathcal{M}$ , and a ciphertext  $\bar{c} \in \mathcal{C}$  such that the two sides of (2.1) are unequal.

Let D be the uniform distribution over  $\mathcal{M}$ , let  $m \in \mathcal{M}$  be arbitrary, let  $\bar{k} \in \mathcal{K}$  be an arbitrary key in the support of Gen (i.e., Gen outputs  $\bar{k}$  with probability > 0), and let  $\bar{c}$  be any ciphertext in the support of  $\mathsf{Enc}_{\bar{k}}(m)$  (keeping in the mind that Enc may be randomized). Now assume without loss of generality that Dec is deterministic (exercise: prove that it's really w.l.o.g., using the correctness of decryption), and define

 $\mathcal{D} = \{ \mathsf{Dec}_k(\bar{c}) : k \in \mathcal{K} \} \subseteq \mathcal{M} \text{ to be the set of all possible decryptions of } \bar{c} \text{ under all possible keys. We must have } |\mathcal{D}| \leq |\mathcal{K}| < |\mathcal{M}|, \text{ which implies that there exists some message } \bar{m} \in \mathcal{M} \setminus \mathcal{D}. \text{ By correctness of decryption, and because } D \text{ is the uniform distribution over } \mathcal{M}, \text{ we have}$ 

$$0 = \Pr_{m \leftarrow D, k \leftarrow \mathsf{Gen}}[m = \bar{m} \mid \mathsf{Enc}_k(m) = \bar{c}] < \Pr_{m \leftarrow D}[m = \bar{m}] = \frac{1}{|\mathcal{M}|},$$

as desired.

Notice that the proof above is "effective" in the sense that  $\bar{m}$  can be computed by explicitly enumerating  $\mathsf{Dec}_k(\bar{c})$  for every  $k \in \mathcal{K}$ . However, if  $\mathcal{K}$  is huge, this task may take too long to be feasible.

# 3 Roadmap to Computational Security

We have seen that perfect secrecy is obtainable, but impractical — the key must be as long as the message. Moreover, the proof of this fact provides an actual attack, which enumerates all possible keys. On the one hand, if the key space is huge (say,  $2^{1000}$ ), the attack may need to run for a long, long time. On the other hand, maybe some schemes fall to faster attacks that don't have to enumerate all keys...

Today (and throughout most of the rest of course), we will be focused on schemes that are meaningfully secure (though not "perfectly" so) as long as the adversary does not employ an absurdly large amount of computation. We have good reason to believe that there are inherent limits on the amount of computation that can be performed in the physical universe — for example, the number of atoms in the observable universe is estimated at around  $10^{80}$ , and we have only about  $10^{10}$  years until our sun runs out of fuel and collapses.

The *computationally secure* schemes we develop will have keys much shorter than the message length, but this is really just the beginning: they will also have absurdly strong (seemingly even paradoxical) security and functionality properties that go far beyond just keeping a message secret.

To get there from here, there are many questions to which we must give precise mathematical answers. here are just a few:

- How do we model "bounded computation"?
- What does it mean for a problem to be *hard* for bounded computation?
- Where might we hope to *find* such hard problems?
- How do we *define* security against computationally bounded adversaries?
- How do we *use* hard problems to design schemes that satisfy our security definitions?

To keep a concrete example in mind, perhaps you have heard the (informal) conjecture that "factoring is hard." We will see how to state this conjecture (and others like it) precisely, and start to use it in cryptography.

# 4 Computational Model

## 4.1 Algorithms

Informally, we all know what an algorithm is: something that you can implement in your favorite programming language, and run on any kind of computer (or network of computers) that you might like. As a formal

mathematical object, an algorithm is a *Turing machine*. In this course we will not need the formal definition too often, and the informal notion will usually suffice.

The running time of an algorithm on an input is the number of "basic" steps it performs before terminating. Basic steps include reading or writing to a location in memory, performing an arithmetic operation (on fixed-size pieces of data), etc. For our purposes, the exact set of basic operations will not be too important, because it is possible to translate between different sets of basic operations with only polynomial slowdown: an algorithm (using one set of basic ops) with running time T can be converted into an algorithm (using another set) with running time  $dT^c$ , for some fixed constants c, d. The running time T(n) of an algorithm as a function of input length n is its maximum running time over all inputs of length n. We say that an algorithm is polynomial-time if its running time  $T(n) = O(n^c)$  for some fixed constant  $c \ge 0$ .

Two aspects of our model of an algorithm require a bit more precision. The first is randomization, that is, we grant algorithms the ability to "flip coins." In the Turing machine view, the machine is augmented with a read-only "random tape," which is initialized with an infinite string of uniformly random and independent bits. Alternatively, we can view the algorithm's randomness as an extra argument  $r \in \{0,1\}^*$ , and we denote the output of  $\mathcal A$  on input x with "coins" r as  $\mathcal A(x;r)$ . (When omitted, r is taken to be a uniformly random string.) Note that  $\mathcal A(x)$  defines a probability distribution over its outputs, where the probability is taken over the random coins. The running time T of a probabilistic algorithm  $\mathcal A$  on an input x is the maximum number of steps performed by  $\mathcal A(x;r)$ , over all random strings r. The notions of running time as a function of the input length n=|x|, and probabilistic polynomial-time (PPT), are defined similarly. Note that a PPT algorithm can only "look at" polynomially many bits of its random string.

In this course, PPT algorithms will be our main model of "efficient computation" for cryptographic schemes. This is primarily to give us a *useful*, *well-behaved abstraction*, which allows us not to get too caught up in the details of implementations. Of course, an algorithm with running time  $n^{100}$  is not very useful in practice, and at times we will take a more precise view of running times.

The second aspect of our model, which we use only for modelling adversaries, is non-uniformity. The idea is that we allow an adversary to have some extra "advice" that depends only on the length of its input. This advice can only increase what the algorithm is capable of computing, so security against non-uniform adversaries is potentially stronger than against only uniform ones. The ability to have advice will also simplify some of our security proofs. More formally, we say that  $\mathcal{A}$  is a non-uniform algorithm if there exists an infinite sequence  $w_1, w_2, \ldots \in \{0, 1\}^*$  so that on input x,  $\mathcal{A}$  is also given the extra argument  $w_{|x|}$ . The notion of PPT for non-uniform algorithms is as above; in particular, note that a non-uniform PPT machine can only "look at" polynomially many bits of its advice string, so without loss of generality we can assume that  $|w_n| = O(n^c)$  for some constant  $c \geq 0$ .

#### 4.2 Asymptotics

We have defined running time as a function of input length n, and we will also frequently do so for probabilities. The length n is frequently called the *security parameter* of a cryptographic scheme, because it (informally) lets us specify "how much" security we want.

We say that a function T(n) is *polynomial*, written  $T(n) = \operatorname{poly}(n)$ , if  $T(n) = O(n^c)$  for some fixed constant c. A nonnegative function  $\nu(n)$  is *negligible*, written  $\nu(n) = \operatorname{negl}(n)$ , if it vanishes faster than the inverse of any polynomial:  $\nu(n) = o(n^{-c})$  for every constant c > 0, or equivalently,  $\lim_{n \to \infty} \nu(n) \cdot n^c = 0$ . We usually apply the concept of a negligible function to quantify the probability of an "extremely rare" event, i.e., one that will effectively "never occur" in a polynomial-time universe. As with polynomial time, the

<sup>&</sup>lt;sup>1</sup>It also makes Turing machines equivalent to "circuit families," which leads to more robust security definitions.

concept of a negligible function is a *useful* and *well-behaved* abstraction, though at times we will need to be more precise in our manipulations of probabilities.

**Question 1.** Is the product of a negligible function and a non-negligible function always negligible, always non-negligible, or neither?

## 5 One-Way Functions

There are many different notions of computational "hardness" in computer science, e.g.: undecidability (exemplified by the Halting Problem), NP-completeness (circuit satisfiability), time/space hierarchies, and others.

In cryptography, we have a few notions of hardness, the most basic of which is *one-wayness*. A *one-way function* (OWF) is often called the "minimal" cryptographic object, because a scheme satisfying almost any interesting computational security notion will imply the existence of a OWF. (This is obviously an informal statement, but a good rule of thumb. We will see some examples later on.) In other words, we can't have (computational) cryptography without one-way functions.

**Definition 5.1.** A function  $f: \{0,1\}^* \to \{0,1\}^*$  is *one-way* if it satisfies the following conditions.

- Easy to compute. There is an efficient algorithm computing f. Formally, there exists a (uniform, deterministic) poly-time algorithm F such that F(x) = f(x) for all  $x \in \{0, 1\}^*$ .
- Hard to invert. An efficient algorithm inverts f on a random input with only negligible probability. Formally, for any non-uniform PPT algorithm  $\mathcal{I}$ , the advantage of  $\mathcal{I}$

$$\mathbf{Adv}_f(\mathcal{I}) := \Pr_{x \leftarrow \{0,1\}^n} \left[ \mathcal{I}(1^n, f(x)) \in f^{-1}(f(x)) \right] = \text{negl}(n)$$

is a negligible function (in the security parameter n).

Some remarks on this definition:

- 1. On  $f^{-1}(f(x))$  versus x: notice that the inverter  $\mathcal{I}$  "wins" if it outputs any preimage of f(x), i.e., any element in the set  $f^{-1}(f(x))$ . (The preimage need not even have length n.) This is mainly to rule out trivial function, such as the function f that just maps every input to 0. If we required the inverter to output the original  $x \in \{0,1\}^n$ , then the best any inverter could do against this f would be to guess randomly, which succeeds with only negligible probability  $2^{-n}$ . But this f does not satisfy our intuitive idea of a function that is "hard to invert."
- 2. The choice of  $\{0,1\}^*$  as the domain and range of f is mainly for syntactic convenience, and it may be replaced by any finite sets, one for each value of the security parameter n. There should also be an efficient algorithm for sampling from the domain, so that the entire inversion experiment can be executed efficiently.
- 3. In contrast to some other notions of hardness you may have seen, this definition is inherently average-case rather than worst-case. A worst-case definition would say that every  $\mathcal{I}$  fails to invert f(x) for some (possibly rare) value of x. Our definition is potentially much stronger: it says that  $\mathcal{I}$  fails to invert f(x) for almost all values of  $x \in \{0,1\}^n$  (for large enough n).

4. Notice that in addition to the value f(x) to be inverted, the input to  $\mathcal{I}$  includes the string  $1^n$ . This is simply a technicality that always allows the inverter to run in time poly(n), even if f(x) has length much less than n. Usually we will omit this extra argument, with the implicit understanding that all algorithms are given the security parameter (in unary), and may run in time polynomial in it.

**Question 2.** Notice that Definition 5.1 allows for  $\mathcal{I}$  to have some non-zero (positive) advantage, though it must be negligible. Consider a stronger definition that requires  $\mathcal{I}$  to have advantage 0 (i.e., it is impossible to find a preimage in polynomial time). Is this definition plausibly satisfiable? Why or why not?

**Question 3.** Let f be a one-way function. Is the function  $g(x) = f(x) \|0$  necessarily one-way?

#### 5.1 Candidates

We have defined the concept of a one-way function, but does such an object exist? First, it can be shown that if a one-way function exists, then  $P \neq NP$ . Since we have no idea how to show that  $P \neq NP$ , proving that a one-way function exists is far beyond our reach. But we have many *candidate* functions that we believe to be one-way. Consider these two examples:

1. Subset-sum: define  $f_{ss}: (\mathbb{Z}_N)^n \times \{0,1\}^n \to (\mathbb{Z}_N)^n \times \mathbb{Z}_N$ , where  $N=2^n$ , as

$$f_{ss}(a_1, \dots, a_n, b_1, \dots, b_n) = (a_1, \dots, a_n, S = \sum_{i: b_i = 1} a_i \mod N).$$

Notice that because the  $a_i$ s are part of the output, inverting f means finding a subset of  $\{1, \ldots, n\}$  (corresponding to those  $b_i$ s equaling 1) that induces the given sum S modulo N.

2. Multiplication: define  $f_{\mathrm{mult}}: \mathbb{N}^2 \to \mathbb{N}$  as  $^2$ 

$$f_{\text{mult}}(x,y) = \begin{cases} 1 & \text{if } x = 1 \lor y = 1 \\ x \cdot y & \text{otherwise.} \end{cases}$$

Are these one-way functions, according to our definition? This is something for you to think about until the next lecture.

<sup>&</sup>lt;sup>2</sup>For security parameter n, we use the domain  $[1, 2^n] \times [1, 2^n]$ . The cases x = 1 and y = 1 are special to rule out the trivial preimages (1, xy) and (xy, 1) for the output  $xy \neq 1$ .

### **Answers**

**Question 1.** Is the product of a negligible function and a non-negligible function always negligible, always non-negligible, or neither?

**Answer.** Neither; it depends on the functions. Consider  $f(n) = 2^{-n}$  (which is negligible) and  $g(n) = 2^n$  (which is non-negligible). Then,  $f(n) \cdot g(n) = 1$ , which is clearly non-negligible. On the other hand, consider h(n) = 1 (which is non-negligible). Then  $f(n) \cdot h(n) = 2^{-n}$ , which is negligible.

**Question 2.** Notice that Definition 5.1 allows for  $\mathcal{I}$  to have some non-zero (positive) advantage, though it must be negligible. Consider a stronger definition that requires  $\mathcal{I}$  to have advantage 0 (i.e., it is impossible to find a preimage in polynomial time). Is this definition plausibly satisfiable? Why or why not?

Answer. It is impossible to satisfy this modified definition. Consider an  $\mathcal{I}$  that ignores the f(x) it is given and simply outputs a uniformly random string from  $\{0,1\}^n$  as its "guess" for a preimage of f(x). This strategy is obviously polynomial time, and it has a nonzero success probability of (at least)  $2^{-n}$ . (There is a  $2^{-n}$  probability of guessing the very same value of x as was used to compute f(x), and there might be other n-bit preimages of f(x) as well, which can only increase the probability of success.) Hence, no function can satisfy the stronger definition. In cryptography, when we can't achieve the "ideal" of making a task completely impossible, we usually do the next-best thing by using a negligible function to mean "effectively zero."

**Question 3.** Let f be a one-way function. Is the function  $g(x) = f(x) \| 0$  necessarily one-way?

**Answer.** Yes, it is. Observe that any g-preimage of g(x) = f(x)||0 is also an f-preimage of f(x), so any potential inverter for g is "effectively" also an inverter for f. Since the latter can have only negligible advantage (because f is one-way), the same goes for the former, and hence g is one-way. In future lectures we will formalize this argument using a *reduction*: for any hypothetical efficient g-inverter  $\mathcal{I}_g$ , we will use it as a "black box" to construct an efficient f-inverter  $\mathcal{I}_f$  such that  $\mathbf{Adv}_f(\mathcal{I}_f) = \mathbf{Adv}_g(\mathcal{I}_g)$ .