CS 341: Algorithms

Charles Shen

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Feel free to email feedback to me at ccshen902@gmail.com.

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1 Selection in deterministic linear time

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In the selection problem, the goal is to determine the i-th smallest element from an unsorted list of n elements in linear time.

"Deterministic" here means that no random numbers are used.

For simplicity, suppose that all elements are distinct, but this is not crucial as a small modification will make it work any list.

One method to choose a pivot is via randomness.

We want a deterministic method to clever choose a pivot to be used to partition a list.

Thus we arrive at the "medians-of-five" algorithm by Blum, Floyd, Pratt, Rivest, and Tarjan (1973).

Idea:

Split the list into $\lfloor \frac{n}{5} \rfloor$ groups of 5 elements, and one additional group of at most 4 elements.

Then sort each group of 5 elements, and find the median of each group of 5. This can be done in constant time, so sorting each group of 5 costs O(n).

Note that that median is the 3rd (center) element in the group of 5. This gives a list of $\lfloor \frac{n}{5} \rfloor$ medians.

Now recursively determine the median of these $\lfloor \frac{n}{5} \rfloor$ medians, which is the element we'll partition around.

We arrive at the algorithm:

```
function SELECT(A, i)
    n \leftarrow |A|
    if n < 60 then
        SORT(A)
        return i-th smallest element
    else
        m \leftarrow \lfloor \frac{n}{5} \rfloor
        divide A up to m groups of 5 elements, with at most one remaining
group of \leq 4 elements
        sort each of the m groups in ascending order
        M \leftarrow \text{array of medians of each group}
        x \leftarrow \text{SELECT}(M)\lceil m/2 \rceil
                                                          ▶ median of all the medians
        k \leftarrow \text{X-PARTITION}(A, x) \Rightarrow \text{partition array A into elements} \leq x \text{ and}
elements > x; returns number of elements on "low side" of the partition
        if i = k then
```

```
return x
      else if i < k then
         return SELECT(A[1..k-1],i)
      else
         return SELECT(A[k+1..n], i-k)
      end if
   end if
end function
```

Let's obtain a lower bound on the number of elements > x, the median of medians.

Here's a picture for n = 37:

smallest	S	S	S	S	*	*	*	*
	S	S	S	S	*	*	*	*
	S	S	S	m	L	L	L	
	*	*	*	L	L	L	L	
largest	*	*	*	L	L	L	L	

m = median of medians

There are $\left\lfloor \frac{n}{5} \right\rfloor$ total columns in which 5 elements appear. Of these, at least $\frac{1}{2}$ (more precisely, $\lceil \lfloor n/5 \rfloor / 2 \rceil$) contain an L. All of these columns, except the one where x appears, contributes 3 to the count of L's; the one where x appears contributes 2.

The conclusion is that $\geq 3 \lceil \lfloor n/5 \rfloor / 2 \rceil - 1$ elements are L's, that is, greater than x. Hence, at most $\leq n - (3 \lceil \lfloor n/5 \rfloor / 2 \rceil - 1)$ elements are $\leq x$.

Then, $n-(3\lceil \lfloor n/5 \rfloor/2 \rceil-1) \leq 7n/10+3$ for all $n \geq 1$. To prove this, see that

$$\lfloor n/5 \rfloor \ge n/5 - 1$$

SO

$$ceilfloorn/5/2 \geq n/10-1/2$$

then

$$3\left\lceil \left\lfloor n/5\right\rfloor /2\right\rceil -1\geq 3n/10-5/2$$

hence

$$n - (3 \lceil |n/5|/2 \rceil - 1) \le 7n/10 + 3$$

We can also claim that at most (7n/10 + 3) elements are $\geq x$ using the same proof above.

So whether we recurse in the smaller elements or larger elements, we can use this bound.

It follows that the time T(n) to select from a list of n elements satisfies the following inequality:

$$T(n) \le T(\lceil n/5 \rceil) + T(r) + dn$$

where $r \leq 7n/10 + 3$, and the dn term soaks up the time needed to do the sorting of 5-element groups, and the partitioning.

Suppose that this recurrence obeys $T(n) \leq cn$ for some c,

$$T(n) \le T(\lfloor n/5 \rfloor) + T(r) + dn$$

$$\le cn/5 + cr + dn$$

$$\le cn/5 + c(7n/10 + 3) + dn$$

$$\le 9cn/10 + 3c + dn$$

and we want this to be less than cn.

Now, $9cn/10 + 3c + dn < cn \iff (c/10 - d)n > 3c$

Now n had better be bigger than 30 here, or else we are in a heap o' trouble. So let's assume $n \ge 60$. Then $(c/10 - d)n \ge (c/10 - d)60$, and if this is to be $\ge 3c$, then we must have $6c - 60d \ge 3c$, and so $c \ge 20d$.

If $n \geq 60$, and $c \geq 20d$, then the induction step indeed works.

To complete the proof, it remains to see that $T(n) \leq cn$ for n = 1, 2, 3, ..., 59. To do this, simply choose

$$c = max(20d, T(1), T(2)/2, ..., T(59)/59)$$

That looks like cheating, but it works.

Now the basis (namely $1 \le n \le 59$) works out just fine, and so does the induction step.

The presence of these large constants (such as c = 20d) suggests that this algorithm is of more theoretical than practical interest.

However, with a bit more work, the number of comparisons can be reduced quite a lot, down to (3 + o(1))n = 3n + o(n).

2 Lower bounds

Can we prove that an algorithm is best possible? Sometimes.

We argued that any comparison-based algorithm for sorting must use $\Omega(n \lg n)$ time. And in reference to the convex hull algorithm, we argued that we couldn't do better than $\Omega(n \lg n)$, because we could use convex hull as a subroutine to create a sorting algorithm; if we could solve convex hull in $o(n \lg n)$ time, then we would end up being able to sort in $o(n \lg n)$ time, a contradiction.

By "lower bounds", we mean a lower bound on the complexity of a *problem*, not an algorithm. Basically we need to prove that *no algorithm*, no matter how complicated or clever, can do better than our bound.

We don't need each algorithm to be "slow" as f(n) on the *same* input, but we do need each algorithm to be "slow" as f(n) on *some* input.

Also, a lower bound doesn't rule out the possibility that some algorithm might be very fast on *some inputs* (but not all).

What do we mean by complexity of a problem?

We say that a function f is a lower bound on the complexity of a problem P if, for every algorithm A to solve P, and every positive integer n, there exists some input I of size n such that A uses at least f(n) steps on input I.

The easiest kind of lower bound is the "information-theoretic bound".

Idea: every algorithm that uses only yes/no questions to decide among k possible alternatives, must ask at least ($\lg k$) questions in the worst case.

Proof. Use an "adversary argument".

The adversary doesn't decide on one of the alternatives ahead of time, but rather it answers the algorithm's questions by always choosing the alternative that maximizes the size of the solution space.

Since the questions are yes/no, the solution space goes down in size by a factor of at most two with each question.

The algorithm cannot answer correctly with certainty unless the solution space is of size 1.

Thus at least $(\lg k)$ questions are needed.

Theorem 2.0.1. Any comparison-based sorting algorithm must use at least $\lg(n!) = \Omega(n \log n)$ questions in the worst-case.

Proof. Use the information-theoretic bound, and observe that the space of possible solutions includes the n! different ways the input could be ordered. \square

3 Adversary Strategy

3.1 Adversary Arguments in General

More generally, we can think of a lower bound proof as a game between the algorithm and an "adversary".

The algorithm is "asking questions" (for example, comparing two elements of an array), while the adversary is answering them (providing the results of the comparison).

The adversary should be thought of as a very powerful, clever being that is answering in such a way as to make your algorithm run as slowly as possible. The adversary cannot "read the algorithm's mind", but it can be prepared for anything the algorithm might do.

Finally, the adversary is not allowed to "cheat"; that is, the adversary cannot answer questions inconsistently. The adversary does not not need to have a particular input in mind at all times when it answers the questions, but when the algorithm is completed, there must be at least one input that matches the answers the adversary gave (otherwise it would have cheated).

The algorithm is trying to run as quickly as possible.

The adversary is trying (through its cleverness) to *force* the algorithm to run slowly.

This interaction proves that the lower-bound obtained by this type of argument applies to *any possible* algorithm from the class under consideration.

Theorem 3.1.1. Every comparison-based algorithm for determining the minimum of a set of n elements must use at least $(\frac{n}{2})$ comparisons.

Proof. Every element must participate in at least one comparison; if not, the not compared element can be chosen (by an adversary) to be the minimum. Each comparison compares 2 elements.

Hence, at least $(\frac{n}{2})$ comparisons must be made.

Theorem 3.1.2. Every comparison-based algorithm for determining the minimum of a set of n elements must use at least (n-1) comparisons.

Proof. To say that a given element, x, is the minimum, implies that (1) every other element has won at least one comparison with another element (not necessarily x). (By "x wins a comparison with y" we mean x > y.) Each comparison produces at most one winner.

Hence at least (n-1) comparisons must be used.

To convert this to an adversary strategy, do the following: for each element, record whether it has won a comparison (W) or not (N).

Initially all elements are labelled with N.

When we compare an N to an N, select one arbitrarily to be the W.

If no values assigned, choose them to make this so.

If values are assigned, decrease the loser (if necessary) to make this so.

When we compare W to N, always make the N the loser and set or adjust values to make sure this is the case. (You might need to decrease the loser's value to ensure this.)

When we compare W to W, use the values already assigned to decide who is the winner.

Now each comparison increases the total number of items labelled W by at most 1.

In order to "know" the minimum, the algorithm must label (n-1) items W, so this proves at least (n-1) comparisons are required.

The values the adversary assigned are the inputs that forced that algorithm to do the (n-1) comparisons.

3.2 Finding Both the Maximum and Minimum of a List of n Numbers

It is possible to compute both the maximum and minimum of a list of n numbers, using $(\frac{3}{2}n-2)$ comparisons if n is even, and $(\frac{3}{2}n-\frac{3}{2})$ comparisons if n is odd.

It can be proven that *no comparison-based method* can correctly determine both the max and min using fewer comparisons in the worst case.

We do this by constructing an adversary argument.

In order for the algorithm to correctly decide that x is the minimum and y is the maximum, it must know that

- 1. every element other than x has won at least one comparison (i.e. x > y if x wins a comparison with y),
- 2. every element other than y has lost at least one comparison

Calling a win W and a loss L, the algorithm must assign (n-1) W's and (n-1) L's. That is, the algorithm must determine (2n-2) "units of information" to always give the correct answer.

We now construct an adversary strategy that will force the algorithm to learn its (2n-2) "units of information" as slowly as possible.

The adversary labels each element of the input as N, W, L, or WL. These labels may change over time.

N signifies that the element has never been compared to any other by the algorithm.

W signifies the element has won at least one comparison.

L signifies the element has lost at least one comparison.

WL signifies the element has won at least one and lost at least one comparison.

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So the the adversary uses the following table.

labels when comparing elements (x, y)	the adversary's response	the new label	unit of information given to the alg.
(N,N)	x > y	(W, L)	2
(W,N) or (WL,N)	x > y	(W, L) or (WL, L)	1
(L,N)	x < y	(L,W)	1
(W,W)	x > y	(W, WL)	1
(L, L)	x > y	(WL, L)	1
(W, L) or (WL, L) or (W, WL)	x > y	no change	0
(WL, WL)	consistent with assigned values	no change	0

The adversary also tentatively assigns values to the elements, which may change over time.

However, they can only change in a fashion consistent with previous answers. That is, an element labelled L may only decrease in value (since it lost all previous comparisons, if it is decreased, it will still lose all of them), and an element labelled W may only increase in value.

An element labelled WL cannot change.

Example. Consider the following sequence of possible questions asked by the algorithm and the adversary's responses:

Algorithm	-	Adve	rsary's r	esponse	s and wor	ksheet		
compares		x1	x2	x3	x4	x5	x6	
(x1, x2)	 	W-20	L-10	N	N	N	N	
(x4, x5)	-	W-20	L-10	N	W-30	L-15	N	
(x1, x4)	-	W-40	L-10	N	WL-30	L-15	N	(
(x3, x6)	-	W-40	L-10	W-11	WL-30	L-15	L-2	
(x2, x5)	-	W-40	WL-10	W-11	WL-30	L-7	L-2	
(x3, x1)	- [WL-40	WL-10	W-50	WL-30	L-7	L-2	
(x2, x4)	-	WL-40	WL-10	W-50	WL-30	L-7	L-2	(
(x5, x6)	-	WL-40	WL-10	W-50	WL-30	WL-7	L-2	

At this point the algorithm knows that x_3 is the maximum, since it is the only element that has never lost a comparison, and x_6 is the minimum, since it is the only element that has never won a comparison.

Note that in step (*), the adversary was forced to reassign the value he had previously assigned to x_1 , since x_1 had to win the comparison against $x_4 = 30$. This is permitted, since x_1 had won every previous comparison and so increasing its value ensures consistency with previous answers.

Also note that step (**) is superfluous, as the algorithm didn't learn any new information.

Theorem 3.2.1. Every comparison-based method for determining both the maximum and minimum of a set of n numbers must use at least $(\frac{3}{2}n-2)$ comparisons (if n even), or $(\frac{3}{2}n-\frac{3}{2})$ comparisons (if n odd), in the worst case.

Proof. Suppose that n is even.

As shown before, the algorithm must learn (2n-2) units of information.

The most it can learn in one comparison is 2 units; when both elements have never participated before in a comparison, labelled by (N, N). This can happen at most $(\frac{n}{2})$ times.

To learn the remaining (n-2) units of info., the algorithm must ask (n-2) questions.

The total number of questions is therefore at least $\frac{n}{2} + n - 2 = \frac{3}{2}n - 2$. The same kind of argument works for n odd.

So we have a lower bound for the problem of finding both the maximum and minimum of a set of n numbers.

So the algorithm is as followed:

If n is even, compare x[1] with x[2] (determining both the maximum and minimum with 1 comparison), x[3] with x[4], etc. We get $(\frac{n}{2})$ maxima and $(\frac{n}{2})$

To find the maximum, use the usual method on the $(\frac{n}{2})$ maxima, which uses

 $(\frac{n}{2}-1)$ comparisons, and the same for the minima. The total cost is $\frac{n}{2} + \frac{n}{2} - 1 + \frac{n}{2} - 1 = \frac{3}{2}n - 2$. A similar method works when n is odd.

Notice that each comparison adds at most one W, and we need to get (n-1)W's before we know the remaining element must be the minimum.

So we need at least (n-1) comparisons.

If the algorithm declares a minimum before getting (n-1) W's, then there are at least two elements without W, so if the algorithm declares one of the minimum, the adversary can truthfully produce the other as the minimum.

3.3 Remarks on Constructing an Adversary Strategy

Lower bound arguments are some of the most difficult and deepest areas in theoretical computer science.

Tips for producing an adversary strategy:

Recall that the lower bound argument can be viewed as a game between you (playing the role of an algorithm) and an adversary.

You want the algorithm to run quickly (for example, by making a small number of queries to the input).

The adversary wants the algorithm to run slowly.

The adversary "wins" (and the lower bound of f(n) is proved) if he/she can force the algorithm to execute f(n) steps.

A proof behaves something like: for all sequences of queries into the input made by the algorithm, there exists a sequence of replies by the adversary, such that the algorithm is forced to execute f(n) steps. So your adversary argument must apply to every possible "move" the algorithm could make.

In constructing its replies, the adversary can do as much or as little bookkeeping as necessary.

There is no requirement that the adversary's computation be efficient or run in any particular time bound.

3.4 Finding the Second Largest

Problem: given a list of n elements, find the 2nd largest.

The naive method (finding the largest, remove it, and then find the largest again) takes 2n-3 comparisons.

Instead, we can do a "tournament" method.

Have the elements in pairs, the larger one (winner) advances; then the 2nd largest (runner-up) is among those defeated by the winner! This is because any other element must have played two other elements superior to it.

This algorithm uses (n-1) comparisons to determine the largest, plus $\lceil \lg n \rceil - 1$ comparisons to determine the runner-up (2nd largest).

The total is $(n + \lceil \lg n \rceil - 2)$ comparisons.

We can prove that the algorithm just given is *optimal*, in the sense that no comparison-based algorithm uses fewer than $(n + \lceil \lg n \rceil - 2)$ comparisons in the worst case.

Note: see Lecture 14 for full proof! (Click here.)

3.5 A Lower Bound for Finding the Median

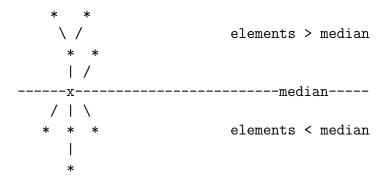
The best algorithm known (in terms of number of comparisons used) uses 3n + o(n) comparisons in the worst case (see Section 1).

The best *lower bound* known for the median-finding problem is 2n + o(n) comparisons.

We will prove a weaker lower bound: $\frac{3}{2}n + O(1)$, where n odd for convenience.

First, note that any comparison-based algorithm for finding the median must use at least (n-1) comparisons. That's because the algorithm must establish the relationship of each element to the median—either by direct comparison, or by comparison with another element whose relationship to the median is known.

The algorithm *must* eventually gain enough knowledge equivalent to the following tree:



Note that if there were an element that didn't participate in a comparison, we couldn't know the median.

For an adversary could change the value of that element from, say, less than the median to greater than the median, changing the median to some other element.

Similarly, if there is some element whose relationship to the median is not known (because, for example, it is greater than some element that is less than the median), we could simply move it from one side of the line above to the other, without changing comparisons. This would also change the median.

Now there are n elements in the tree above, so there are (n-1) edges, and hence (n-1) comparisons.

We now improve this to $\frac{3}{2}n + O(1)$.

First, we call a comparison involving x crucial, say x against y, if it is the first comparison where

- 1. x < y and $y \le \text{median OR}$ if
- 2. x > y and $y \ge$ median

So the algorithm must perform at least (n-1) crucial comparisons in order to determine the median.

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We now give an adversary strategy to *force* the algorithm to perform (n-1)/2 non-crucial comparisons (since we've proved earlier that the algorithm requires (n-1) crucial comparisons.

Idea:

First, the adversary picks some value (not some element) to be the median.

Then it assigns a label in $\{N, L, S\}$ to each element, and also values, as appropriate.

Initially each element is assigned an N ("Never participated in any comparison"). L means Larger than the median and S means Smaller than the median. The adversary follows the following strategy:

Algorithm compares	Adversary responds
(N, N)	assign one element to be larger than the median, one smaller; result is (L, S)
(S,N) or (N,S)	assign the N-element to be larger than the median; result is (S, L) or (L, S)
(L,N) or (N,L)	assign the N-element to be smaller than the median; result is (L, S) or (S, L)
(S, L) or (L, S) or (S, S) or (L, L)	consistent with previously assigned values

Note 1: this strategy continues until ((n-1)/2) S's (or ((n-1)/2) L's) have been assigned. If at some point ((n-1)/2) S's are assigned, then the adversary assigns the remaining elements to be greater than the median, except for one, which is the median. A similar thing is done if ((n-1)/2) L's have been assigned.

Note 2: the last element assigned is always the median.

I claim that this strategy will always force the algorithm to perform ((n-1)/2) non-crucial comparisons.

For any time an N-element is compared, a non-crucial comparison is done (except at the very end, when a crucial comparison may be done with the median itself).

The least number of comparisons with N-elements that can be done is ((n-1)/2).

The total number of comparisons is therefore n-1+(n-1)/2=(3/2)(n-1).

Example. The adversary strategy when n = 7. The adversary arbitrarily chooses the median to be 43. Then

Algorithm Adversary's strategy (label:value) compares x1 x2 x3 x4 x5 x6 x7

	N	N	N	N	N	N	N	
(x1, x3)	L:50	N	S:31	N	N	N	N	
(x3, x4)	L:50	N	S:31	L:47	N	N	N	
(x6, x7)	L:50	N	S:31	L:47	N	L:60	S:20	
	abbighou, by davoibary abbighb one							
(x2, x5)	a crud	cial	S:31 compari	son	S:15	L:60	S:20	
-	(x2, x1) a crucial comparison							
(x2, x3) a crucial comparison								
(x2, x4)	(x2, x4) a crucial comparison							
(x2, x6)	x6) a crucial comparison							
(x2, x7)	a crud	cial	compari	son				

At this point the algorithm concludes that x2 is the median.

The first three comparisons were non-crucial; only the last six are crucial.

The best lower bound known for the median: Dor and Zwick, Median selection requires $(2 + \epsilon)n$ comparisons.

Lower Bound on Average-Case Sorting 3.6

Theorem 3.6.1. Any comparison-based algorithm must use $\Omega(n \lg n)$ comparisons on average.

Lemma 3.6.2. In any binary tree T where each node has 0 or 2 children,

$$\sum_{\ell \text{ a leaf of } T} 2^{-\operatorname{depth}(T,\ell)} = 1$$

where $depth(T, \ell)$ is the length of the path from ℓ to the root.

Proof. Prove by induction on the height of T.

Easy to see the lemma holds when T consists of a single node.

Assume it holds for all tree of height $\langle k \rangle$; prove it for trees of height k.

Then a tree of height k consists of a root and two subtrees, T' and T''.

Then the lemma applies to each of these subtrees recursively, so we know that

$$\sum_{\ell \text{ a leaf of } T'} 2^{-\operatorname{depth}(T',\ell)} = 1$$

and

$$\sum_{\ell \text{ a leaf of } T''} 2^{-\operatorname{depth}(T'',\ell)} = 1$$

Then we have

$$\sum_{\ell \text{ a leaf of } T'} 2^{-\operatorname{depth}(T,\ell)} = \frac{1}{2}$$

since the depth of a leaf in T goes up by 1 when considered as a leaf of T. The same result holds for T''.

Adding these together yields the result.

Lemma 3.6.3. Suppose $a, b \in \mathbb{R}$ and a, b > 0. If $a \neq b$, then we have

$$\lg a - \lg b > 2\lg((a+b)/2)$$

Proof. If $a \neq b$, then

$$(\sqrt{a} - \sqrt{b})^2 > 0$$

$$a - 2\sqrt{ab} + b > 0$$

$$(a+b)/2 > \sqrt{ab}$$

$$\lg((a+b)/2) > \frac{1}{2}\lg(ab)$$

$$2\lg((a+b)/2) > \lg a + \lg b$$

$$-2\lg((a+b)/2) < -\lg a - \lg b$$

as required.

Lemma 3.6.4. Suppose $x_1, x_2, \ldots, x_k \in \mathbb{R}$ and $x_1, x_2, \ldots, x_k > 0$ such that $\sum_{1 \leq i \leq k} x_i = 1$. Then

$$\sum_{1 \le i \le k} (-\lg x_i) \ge k \lg k$$

Proof. If all the $x_i = \frac{1}{k}$, then the inequality clearly holds - actually then it is an equality!

Suppose we claim this sum actually achieves its minimum at this point.

For if the minimum occurred at some *other* point, then two of the x_i must be different, say a and b.

Then by Lemma 3.6.3, we could replace $(-\lg a) + (-\lg b)$ with two copies of $-\lg((a+b)/2)$ and get a smaller sum, a contradiction.

Proof. Consider the comparison tree used by the algorithm.

There are at least k = n! leaves, since there are n! possible ways to order n numbers.

There are also at most this number of leaves, since otherwise two leaves would be labelled with the same ordering, and then, by tracing up to the lowest common ancestor, we would get to a node where there are two different answers, implying the orderings cannot be the same.

So there are exactly n! leaves.

By Lemma 3.6.2,

$$\sum_{\ell \text{ a leaf of } T} 2^{-\operatorname{depth}(\ell)} = 1$$

Then by Lemma 3.6.4, with $x_i = 2^{-\operatorname{depth}(\ell)}$, we get

$$\sum_{\substack{\ell \text{ a leaf of } T}} \operatorname{depth}(\ell) \ge n! \lg(n!)$$

It follows that

$$\frac{1}{n!} \sum_{\ell \text{ a leaf of } T} \operatorname{depth}(\ell) \ge \lg(n!)$$

but the expression on the left is just the average depth of each leaf, i.e., the average number of comparisons used (averaged over all possible arrangements of the input).

Hence the average number of comparisons used is $\Omega(n \lg n)$.

4 Graphs and Graph Algorithms

A graph consists of a bunch of points, called vertices (or nodes), and a bunch of connections between pairs of points, called edges (or arcs).

So G = (V, E), where V is a set of vertices and E is a subset of $V \times V$.

Where |V| is the number of vertices and |E| is the number of edges in G.

Graphs can be directed or undirected; in a directed graph, edges have a direction, and you can only traverse the graph by going in that direction. In an undirected graph, edges can be traversed in any direction.

Sometimes we allow graphs to have edges from a vertex to itself; such an edge is called a "loop" or a "self-loop" and the graph is no longer *simple*.

Most graphs in this course won't have such edges, we will allow self-loops for directed graphs and only sometimes for undirected graphs. We will rarely have two or more different edges between the same two vertices.

Note: in any graph, $|E| = O(|V|^2)$, and if the graph is connected, $|E| = \Omega(|V|)$.

A *path* is a series of vertices where each consecutive vertex is connected by an edge to the previous vertex.

An undirected graph is *connected* if there is a path from every vertex to every other vertex.

The *length* of a path is the number of edges in the path.

A path is a *cycle* if the first vertex is the same as the last vertex. Usually in a path we demand that there are no repeated vertices (with the exception in a cycle).

A walk (resp., closed walk) is a path (resp., cycle) where repeated vertices and edges are allowed.

A graph is *acyclic* if there are no cycles.

A tree is a connected acyclic graph.

An undirected graph is a tree if every pair of vertices are connected by a *unique* simple path.

Trees can be rooted (have a distinguishable vertex called the root) or unrooted.

Two ways to store a graph:

1. Adjacency Matrix

We represent a graph by a $|V| \times |V|$ matrix with a 1 in row i and column

j is (i, j) is an edge, and 0 otherwise.

2. Adjacency List

We represent a graph by an array Adj of |V| lists, one list for each vertex. For each vertex Adj[u] contains all vertices (or pointers to the vertices) v such that uv is an edge.

Advantages and disadvantages of each representation: $Adjacency\ Matrix$:

- easy to manipulate
- can check if uv is an edge in O(1) time
- but storage is always $\Omega(|V|^2)$ —not good for sparse graphs
- time to add an edge:
 - existing vertex and new edge: O(1) time
 - existing vertex and existing edge: O(1) time, can check in O(1) time
 - new vertex and new edge: $O(|V|^2)$ time, create a new, bigger matrix

Adjacency List:

- harder to manipulate
- checking if uv is an edge takes O(|V|)time
- but storage is O(|V| + |E|)
- time to add an edge:
 - existing vertex and new edge: O(|V|) time
 - existing vertex and existing edge: O(|V|) time
 - new vertex and new edge: O(|V|) time

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Algorithms for minimum spanning tree:

Kruskal: many different trees that are eventually joined up to form one tree

Prim: one tree that grows by sequentially adding edges

Neither Kruskal nor Prim work for directed graphs.

That problem is called the "minimum cost arborescence" problem; it can be solved in $min(n^2, m \lg n)$ time.

4.1 Kruskal's Algorithm for Minimum Spanning Tree

We consider each edge in turn, starting with the lowest-weight edge and continuing in ascending order by weight. We only add the edge if it doesn't create a cycle.

To implement Kruskal's algorithm, we need to consider the edges in turn, ordered in ascending order by weight; which we can do by sorting the edges. Then when we pick an edge uv to consider, we have to decide if they are in different trees, or the same tree. This is not so easy! But it is an example of the "disjoint-set data structure problem."

In the disjoint-set data structure problem, we have three operations:

MAKE-SET: create a set containing a single vertex

FIND: given a vertex, find the "name" of the set it is in (sets are disjoint so only one name is possible for each vertex)

UNION: union together two sets, possibly giving the resulting set a new name (destructive)

Due to the two main operations, this is also called the "union-find" problem.

It is possible to do m MAKE-SET, UNION and FIND operations,n of which are MAKE-SET, in $O(m \ \alpha(n))$ time.

 $\alpha(n)$ is an extremely slow-growing function called the "inverse Ackerman" function; the function grows so slowly that $\alpha(n) \leq 4$ for all values of n that one would ever encounter in real life.

How Kruskal's algorithm may be implemented:

```
function KRUSKAL(G, w) 
ightharpoonup G = (V, E) and w is the weight function A \leftarrow \text{empty set} for each v in V do MAKE-SET(v) end for sort E into increasing order of weight for each edge uv in E, in increasing order do if \text{FIND}(u) \neq \text{FIND}(v) then A \leftarrow A \cup \{uv\} UNION(u, v)
```

end if end for return A end function

MAKE-SET(v) creates a new set whose only element is v.

FIND(u) returns the name of the set that u is contained in.

UNION(u,v) unions together the trees containing u and v and gives the result a name (which can be found with FIND).

Sets are stored as linked lists of vertices.

Each element in the list has a pointer to its next element, and a pointer to the head of the list.

There is also a tail pointer to the end of the list, and the list maintains how many elements it contains.

The "name" is the first vertex in the list.

To do MAKE-SET, we create a linked list of a single element, and the num field set to 1. This costs O(1).

To do a FIND, we follow the pointer to the head of the list and return the element there. This costs O(1).

To do a UNION(u,v), we follow the pointer of u to its head, and the pointer of v to its head.

There we find the number of elements in each list.

We can do this via "weighted-union heuristic."

If (say) u is in the smaller list, we link that list to the end of the bigger list (v's list).

Then traverse u's list, moving each head pointer to point to the head of v's list.

Finally, we update the num field of v's list.

This costs O(t), where t is the number of vertices of the smaller list.

Because we always update the head pointers of the elements of the smaller list, each UNION at least doubles the size of the smaller list.

So an element's name gets updated at most O(logn) times.

Thus, over the course of the entire algorithm, we see that (all) n UNIONs take at most $O(n \lg n)$ time.

Theorem 4.1.1. Kruskal's algorithm takes $O(|E| \lg |E|) = O(|E| \lg |V|)$ time.

Proof. First, note that the algorithm does indeed produce a spanning tree.

The output cannot have a cycle, because if it did, the edge added that forms the cycle would join together two vertices in the same tree.

The output must be connected, for if it weren't, when we considered the edge that connected two components, we would have added it.

Next, we need to argue that the spanning tree T produced by Kruskal's algorithm is a minimum spanning tree.

Assume it isn't. Among all minimum spanning trees for G, the input, let T' be the one with the largest number of edges in T.

If T' = T, we're done.

Otherwise, among all the edges in T-T', let e be the one that is added earliest by Kruskal's algorithm when we run it on G.

Because T' is a spanning tree, $T' \cup \{e\}$ has a cycle C.

T cannot contain all the edges of this cycle (otherwise it would not be a tree) so C contains an edge f that is not in T.

Now consider $T'' = (T' \cup \{e\}) - \{f\}$. T" is also a spanning tree; since T' was minimum, $weight(e) \ge weight(f)$.

If weight(f) < weight(e), then the algorithm would have considered the edge f before it would have considered the edge e.

Adding f wouldn't create a cycle (e was the first edge not in T'), so it would have been added.

But it wasn't; e was.

So $weight(f) \ge weight(e)$.

Thus weight(e) = weight(f).

So T'' is also a minimum spanning tree.

But T'' has one more edge in common with T than T' does, contradicting the choice of T'.

4.2 Prim's Algorithm

It is a greedy algorithm where, at each step, we choose a previously unconnected vertex that becomes connected by a lowest-weight edge.

We choose an initial vertex arbitrarily and put it in a set V'.

We also have a set E' of edges which initially empty.

At every step, we choose an edge uv of minimal weight with u in V' and v not in V' (break ties arbitrarily).

We then add v to V' and uv to E'.