

CS 473: Algorithms

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Fall 2016

Review session

Lecture 99

September 30, 2016

What we saw so far...

Fast Fourier Transform (FFT).

Dynamic Programming

- String algorithms.
- Graph algorithms: shortest path, independent set, dominating set, etc.

Randomized Algorithms

- Quick sort,
- High probability analysis: Markov, Chebyshev, and Chernoff inequalities

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Randomized Algorithms

- Quick sort,
- High probability analysis: Markov, Chebyshev, and Chernoff inequalities
- Hashing, Fingerprinting

Part I

FFT

What is Fast Fourier Transform

Definition

Given vector $\mathbf{a} = (a_0, a_1, \dots, a_{n-1})$ the *Discrete Fourier Transform* (DFT) of \mathbf{a} is the vector $\mathbf{a}' = (a'_0, a'_1, \dots, a'_{n-1})$ where $a'_j = a(\omega_n^j)$ for $0 \leq j < n$.

\mathbf{a}' is a sample representation of polynomial with coefficient representation \mathbf{a} at n 'th roots of unity.

We have shown that \mathbf{a}' can be computed from \mathbf{a} in $O(n \log n)$ time. This divide and conquer *algorithm* is called the *Fast Fourier Transform* (FFT).

Why FFT? Convolution and Polynomial Multiplication

Convolution

Convolution of vectors $\mathbf{a} = (a_0, a_1, \dots, a_{n-1})$ and $\mathbf{b} = (b_0, b_1, \dots, b_{n-1})$ is a vector $\mathbf{c} = (c_0, c_1, \dots, c_{2n-2})$, where

$$c_k = \sum_{i,j: i+j=k} a_i \cdot b_j$$

(Handwritten red annotations: c_k is circled; $a_i \cdot b_j$ is circled; the term $(x^i \cdot x^j)$ is written in red with x^k below it, indicating the polynomial multiplication context.)

Why FFT? Convolution and Polynomial Multiplication

Convolution

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Polynomial Multiplication

If vectors \mathbf{a} and \mathbf{b} are coefficients of two $n - 1$ degree polynomials, (abusing notation) $\mathbf{a}(\mathbf{x}) = \sum_{i=0}^{n-1} a_i x^i$, $\mathbf{b}(\mathbf{x}) = \sum_{i=0}^{n-1} b_i x^i$ then \mathbf{c} is the coefficient vector of the product polynomial $\mathbf{a}(\mathbf{x}) * \mathbf{b}(\mathbf{x})$.

Why FFT? Convolution and Polynomial Multiplication

Convolution

Given vectors $\mathbf{a} = (a_0, a_1, \dots, a_{n-1})$ and $\mathbf{b} = (b_0, b_1, \dots, b_{n-1})$ find its convolution vector $\mathbf{c} = (c_0, c_1, \dots, c_{2n-2})$.

- 1 Compute values of \mathbf{P}_a and \mathbf{P}_b at the $2n$ th roots of unity, to get their sample representation \mathbf{a}' and \mathbf{b}' .
- 2 Compute sample representation $\mathbf{c}' = \mathbf{a}' \mathbf{b}'_0, \dots, \mathbf{a}'_{2n-2} \mathbf{b}'_{2n-2}$ of product $\mathbf{c} = \mathbf{a} \cdot \mathbf{b}$
- 3 Compute \mathbf{c} from \mathbf{c}' using inverse Fourier transform.

- Step 1 takes $O(n \log n)$ using two FFTs
- Step 2 takes $O(n)$ time
- Step 3 takes $O(n \log n)$ using one FFT

$$C(\omega) = A(\omega) \cdot B(\omega)$$

Problem

Suppose we are given a bit string $B[1..n]$. A triple of distinct indices $1 \leq i < j < k \leq n$ is called a well-spaced triple in B if

$B[i] = B[j] = B[k] = 1$ and $k - j = j - i = d$

- (a) Describe a brute-force algorithm to determine whether B has a well-spaced triple in $O(n^2)$ time.

$B = 010111011$
1 2 3 4 5 6 7 8 9
 / / /
 / / /
For $i = 1 \dots n-3$
 For $j = i+1 \dots n-2$
 For $k = j+1 \dots n$
 If $B[i] = B[j] = B[k] = 1$ and $k - j = j - i$ then return true
 End For
 End For
End For

i, d
2 2
4 1
4 2
 $k = i + 2d \leq n$
 $d \leq n/2$

Application of FFT

Suppose we are given a bit string $B[1..n]$. A triple of distinct indices $1 \leq i < j < k \leq n$ is called a well-spaced triple in B if

$B[i] = B[j] = B[k] = 1$ and $k - j = j - i + a \rightarrow \textcircled{0}$

(b) Describe an algorithm to determine whether B has a well-spaced triple in $O(n \log n)$ time.

$$B[i]B[j]B[k] = 1 \quad \text{if} \quad \textcircled{1}$$

$$\textcircled{2} \quad 2j - i - k + a = 0 \Leftrightarrow 2j - i - k = -a$$

$$P(x) = \sum_{i,j,k} (B[i]B[j]B[k]) \frac{x^j x^{-i} x^{-k}}{x^{-a(2n)}} x^{2n}$$

i, j, k \Rightarrow x

$$P_1(x) = \sum_{j=1}^n B[j] x^{2j}$$

$$P_2(x) = \left(\sum_{i=1}^n B[i] x^{-i} \right) x x^n$$

$$P_3(x) = \left(\sum_{k=1}^n B[k] x^{-k} \right) x x^n$$

Application of FFT

Suppose we are given a bit string $B[1..n]$. A triple of distinct indices $1 \leq i < j < k \leq n$ is called a well-spaced triple in B if $B[i] = B[j] = B[k] = 1$ and $k - j = j - i$.

- (c) Describe an algorithm to determine the number of well-spaced triples in B in $O(n \log n)$ time.

coefficient of x^{2n-a}

Part II

Dynamic Programming

Recursion

Reduction:

Reduce one problem to another

Recursion

A special case of reduction

- 1 reduce problem to a *smaller* instance of *itself*
- 2 self-reduction

- 1 Problem instance of size **n** is reduced to one or more instances of size **n - 1** or less.
- 2 For termination, problem instances of small size are solved by some other method as **base cases**.

What is Dynamic Programming?

Every recursion can be memoized. Automatic memoization does not help us understand whether the resulting algorithm is efficient or not.

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Dynamic Programming:

A recursion that when memoized leads to an *efficient* algorithm.

Edit Distance

Definition

Edit distance between two words **X** and **Y** is the number of letter insertions, letter deletions and letter substitutions required to obtain **Y** from **X**.

Example

The edit distance between FOOD and MONEY is at most **4**:

FOOD → MOOD → MONOD → MONED → MONEY

Edit Distance: Alternate View

Alignment

Place words one on top of the other, with gaps in the first word indicating insertions, and gaps in the second word indicating deletions.

F	O	O		D
M	O	N	E	Y

Edit Distance: Alternate View

Alignment

Place words one on top of the other, with gaps in the first word indicating insertions, and gaps in the second word indicating deletions.

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M	O	N	E	Y

Formally, an **alignment** is a set **M** of pairs **(i, j)** such that each index appears at most once, and there is no “crossing”: **i < i'** and **i** is matched to **j** implies **i'** is matched to **j' > j**. In the above example, this is **M = {(1, 1), (2, 2), (3, 3), (4, 5)}**.

Edit Distance: Alternate View

Alignment

Place words one on top of the other, with gaps in the first word indicating insertions, and gaps in the second word indicating deletions.

F	O	O		D
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Formally, an **alignment** is a set M of pairs (i, j) such that each index appears at most once, and there is no “crossing”: $i < i'$ and i is matched to j implies i' is matched to $j' > j$. In the above example, this is $M = \{(1, 1), (2, 2), (3, 3), (4, 5)\}$. Cost of an alignment is the number of mismatched columns plus number of unmatched indices in both strings.

Edit Distance Problem

Problem

Given two words, find the edit distance between them, i.e., an alignment of smallest cost.

Edit Distance

Basic observation

Let $\mathbf{X} = \alpha\mathbf{x}$ and $\mathbf{Y} = \beta\mathbf{y}$

α, β : strings. \mathbf{x} and \mathbf{y} single characters.

Possible alignments between \mathbf{X} and \mathbf{Y}

α	\mathbf{x}
β	\mathbf{y}

or

α	\mathbf{x}
$\beta\mathbf{y}$	

or

$\alpha\mathbf{x}$	
β	\mathbf{y}

Observation

Prefixes must have optimal alignment!

Edit Distance

Basic observation

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α	\mathbf{x}
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or

α	\mathbf{x}
$\beta\mathbf{y}$	

or

$\alpha\mathbf{x}$	
β	\mathbf{y}

Observation

Prefixes must have optimal alignment!

$$\text{EDIST}(\mathbf{X}, \mathbf{Y}) = \min \begin{cases} \text{EDIST}(\alpha, \beta) + [\mathbf{x} \neq \mathbf{y}] \\ 1 + \text{EDIST}(\alpha, \mathbf{Y}) \\ 1 + \text{EDIST}(\mathbf{X}, \beta) \end{cases}$$

Recursive Algorithm

Assume **X** is stored in array **A[1..m]** and **Y** is stored in **B[1..n]**

```
EDIST(A[1..i], B[1..j])
```

```
  If (i = 0) return j
```

```
  If (j = 0) return i
```

```
  m1 = 1 + EDIST(A[1..(i - 1)], B[1..j])
```

```
  m2 = 1 + EDIST(A[1..i], B[1..(j - 1)])
```

```
  If (A[m] = B[n]) then
```

```
    m3 = EDIST(A[1..(i - 1)], B[1..(j - 1)])
```

```
  Else
```

```
    m3 = 1 + EDIST(A[1..(i - 1)], B[1..(j - 1)])
```

```
  return min(m1, m2, m3)
```

Call **EDIST(A[1..m], B[1..n])**

Memoizing the Recursive Algorithm

```
int M[0..m][0..n]
```

```
Initialize all entries of M[i][j] to  $\infty$   
return EDIST(A[1..m], B[1..n])
```

```
EDIST(A[1..i], B[1..j])
```

```
  If (M[i][j] <  $\infty$ ) return M[i][j]    (* return stored value *)
```

```
  If (i = 0)
```

```
    M[i][j] = j
```

```
  ElseIf (j = 0)
```

```
    M[i][j] = i
```

```
  Else
```

```
    m1 = 1 + EDIST(A[1..(i - 1)], B[1..j])
```

```
    m2 = 1 + EDIST(A[1..i], B[1..(j - 1)])
```

```
    If (A[i] = B[j]) m3 = EDIST(A[1..(i - 1)], B[1..(j - 1)])
```

```
    Else m3 = 1 + EDIST(A[1..(i - 1)], B[1..(j - 1)])
```

```
    M[i][j] = min(m1, m2, m3)
```

```
  return M[i][j]
```

Removing Recursion to obtain Iterative Algorithm

```
EDIST(A[1..m], B[1..n])  
  int M[0..m][0..n]  
  for i = 0 to m do M[i, 0] = i  
  for j = 0 to n do M[0, j] = j  
  
  for i = 1 to m do  
    for j = 1 to n do  
      
$$M[i][j] = \min \begin{cases} [x_i \neq y_j] + M[i-1][j-1], \\ 1 + M[i-1][j], \\ 1 + M[i][j-1] \end{cases}$$

```

Removing Recursion to obtain Iterative Algorithm

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  int M[0..m][0..n]  
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Analysis

Removing Recursion to obtain Iterative Algorithm

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EDIST(A[1..m], B[1..n])  
  int M[0..m][0..n]  
  for i = 0 to m do M[i, 0] = i  
  for j = 0 to n do M[0, j] = j  
  
  for i = 1 to m do  
    for j = 1 to n do  
      
$$M[i][j] = \min \begin{cases} [x_i \neq y_j] + M[i-1][j-1], \\ 1 + M[i-1][j], \\ 1 + M[i][j-1] \end{cases}$$

```

Analysis

- 1 Running time is $O(mn)$.
- 2 Space used is $O(mn)$.

Matrix and DAG of Computation

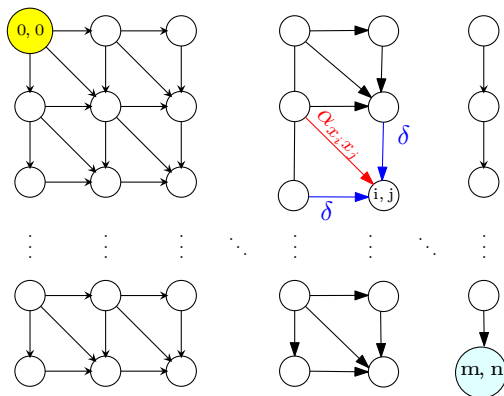


Figure : Iterative algorithm in previous slide computes values in row order.

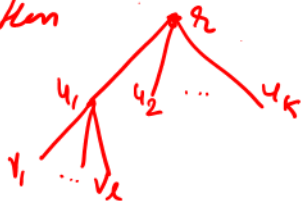
Problem

Given a graph $G = (V, E)$ a matching is a set of edges $M \subset E$ such that no two edges in M share an end point. Describe an efficient algorithm that given a tree $T = (V, E)$ and non-negative weights $w : E \rightarrow \mathbb{R}^+$ finds a maximum weight matching in T .

$M(u) = \text{Weight of the max matching in subtree of } u.$

if u not matched then

$$M(u) = \sum_{v \in C(u)} M(v)$$



if u matched with u_1 then

$$M(u) = w(u, u_1) + \sum_{v \in C(u)} M(v) + \sum_{\substack{v \in C(u) \\ v \neq u_1}} M(v)$$

$$M(x) = \max \left\{ \begin{array}{l} \sum_{u \in C(x)} M(u) \quad x \text{ is internal} \\ \forall u \in C(x) : \\ w(x, u) + \sum_{\substack{u' \neq u \\ u' \in C(x)}} M(u') \\ + \sum_{v \in C(u)} M(v) \end{array} \right.$$

Base case:

$$u \in \text{Leaf}, M(u) = 0$$

Call $M(x)$.

Dijkstra's Algorithm

Initialize for each node v , $\text{dist}(s, v) = \infty$

Initialize $S = \{s\}$, $\text{dist}(s, s) = 0$

for $i = 1$ to $|V|$ **do**

 Let v be such that $\text{dist}(s, v) = \min_{u \in V-S} \text{dist}(s, u)$

$S = S \cup \{v\}$

for each u in $\text{Adj}(v)$ **do**

$\text{dist}(s, u) = \min(\text{dist}(s, u), \text{dist}(s, v) + \ell(v, u))$

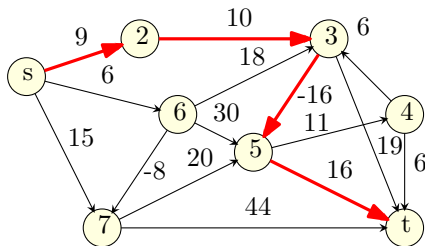
- ① Using Fibonacci heaps. Running time: $O(m + n \log n)$.
- ② Can compute shortest path tree.

Single-Source Shortest Paths with Negative Edge Lengths

Single-Source Shortest Path Problems

Input: A *directed* graph $G = (V, E)$ with arbitrary (including negative) edge lengths. For edge $e = (u, v)$, $\ell(e) = \ell(u, v)$ is its length.

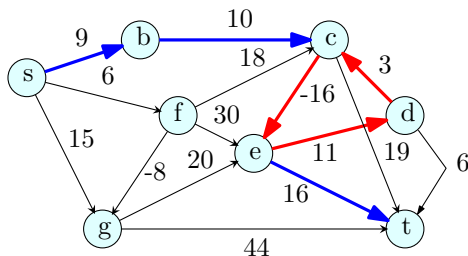
- Given nodes s, t find shortest path from s to t .
- Given node s find shortest path from s to all other nodes.



Negative Length Cycles

Definition

A cycle **C** is a negative length cycle if the sum of the edge lengths of **C** is negative.



Dijkstra's algorithm does not work with negative edges.

Shortest Paths and Recursion

- 1 Compute the shortest path distance from **s** to **t** recursively?
- 2 What are the smaller sub-problems?

Shortest Paths and Recursion

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- 2 What are the smaller sub-problems?

Lemma

Let **G** be a directed graph with arbitrary edge lengths. If $s = v_0 \rightarrow v_1 \rightarrow v_2 \rightarrow \dots \rightarrow v_k$ is a shortest path from **s** to v_k then for $1 \leq i < k$:

- 1 $s = v_0 \rightarrow v_1 \rightarrow v_2 \dots \rightarrow v_i$ is a shortest path from **s** to v_i



Shortest Paths and Recursion

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Lemma

Let **G** be a directed graph with arbitrary edge lengths. If $\mathbf{s} = \mathbf{v}_0 \rightarrow \mathbf{v}_1 \rightarrow \mathbf{v}_2 \rightarrow \dots \rightarrow \mathbf{v}_k$ is a shortest path from **s** to \mathbf{v}_k then for $1 \leq i < k$:

- 1 $\mathbf{s} = \mathbf{v}_0 \rightarrow \mathbf{v}_1 \rightarrow \mathbf{v}_2 \dots \rightarrow \mathbf{v}_i$ is a shortest path from **s** to \mathbf{v}_i

Sub-problem idea: paths of fewer hops/edges

Hop-based Recursion: Bellman-Ford Algorithm

Single-source problem: fix source s .

Assume that all nodes can be reached by s in G . (Remove nodes unreachable from s).

$d(v, k)$: shortest walk length from s to v using at most k edges.

Hop-based Recursion: Bellman-Ford Algorithm

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Recursion for $d(v, k)$:

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$d(v, k)$: shortest walk length from s to v using at most k edges.

Recursion for $d(v, k)$:

$$d(v, k) = \min \begin{cases} \min_{u \in v} (d(u, k-1) + \ell(u, v)). \\ d(v, k-1) \end{cases}$$

Base case: $d(s, 0) = 0$ and $d(v, 0) = \infty$ for all $v \neq s$.

A Basic Lemma

Lemma

Assume s can reach all nodes in $G = (V, E)$. Then,

- ① There is a negative length cycle in G iff $d(v, n) < d(v, n - 1)$ for some node $v \in V$.
- ② If there is no negative length cycle in G then $\text{dist}(s, v) = d(v, n - 1)$ for all $v \in V$.

Bellman-Ford Algorithm

```
for each  $u \in V$  do
     $d(u, 0) \leftarrow \infty$ 
 $d(s, 0) \leftarrow 0$ 

for  $k = 1$  to  $n$  do
    for each  $v \in V$  do
         $d(v, k) \leftarrow d(v, k - 1)$ 
        for each edge  $(u, v) \in \text{In}(v)$  do
             $d(v, k) = \min\{d(v, k), d(u, k - 1) + \ell(u, v)\}$ 

for each  $v \in V$  do
     $\text{dist}(s, v) \leftarrow d(v, n - 1)$ 
    If  $d(v, n) < d(v, n - 1)$ 
        Return ‘‘Negative Cycle in  $G$ ’’
```

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Running time:

Bellman-Ford Algorithm

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Running time: $O(mn)$

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```

Running time: $O(mn)$ Space:

Bellman-Ford Algorithm

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Running time: $O(mn)$ Space: $O(n^2)$

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for each  $v \in V$  do
     $\text{dist}(s, v) \leftarrow d(v, n - 1)$ 
    If  $d(v, n) < d(v, n - 1)$ 
        Return ‘‘Negative Cycle in  $G$ ’’
```

Running time: $O(mn)$ Space: $O(n^2)$

Space can be reduced to $O(m + n)$.

Bellman-Ford with Space Saving

```
for each  $u \in V$  do
     $d(u) \leftarrow \infty$ 
 $d(s) \leftarrow 0$ 

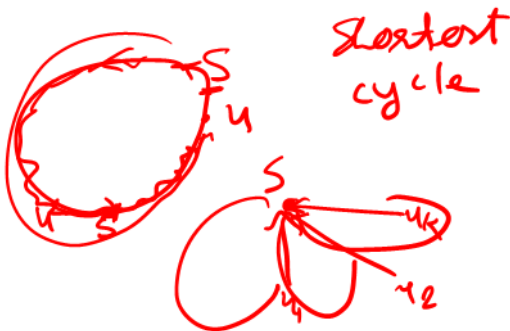
for  $k = 1$  to  $n - 1$  do
    for each  $v \in V$  do
        for each edge  $(u, v) \in \text{In}(v)$  do
             $d(v) = \min\{d(v), d(u) + \ell(u, v)\}$ 

(* One more iteration to check if distances change *)
for each  $v \in V$  do
    for each edge  $(u, v) \in \text{In}(v)$  do
        if  $(d(v) > d(u) + \ell(u, v))$ 
            Output "Negative Cycle"

for each  $v \in V$  do
     $\text{dist}(s, v) \leftarrow d(v)$ 
```

Problem

Given a directed graph $G = (V, E)$ with non-negative edge lengths $\ell : E \rightarrow \mathbf{R}^+$, describe an algorithm that finds the shortest cycle in G that contains a specific node s .



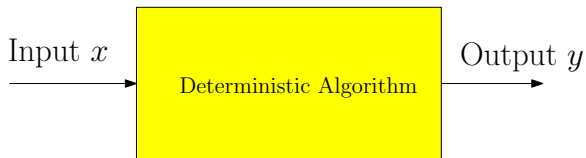
Problem

Given a directed graph $G = (V, E)$ with non-negative edge lengths $\ell : E \rightarrow \mathbf{R}^+$. Describe an algorithm to find the shortest cycle containing s with at most k edges.

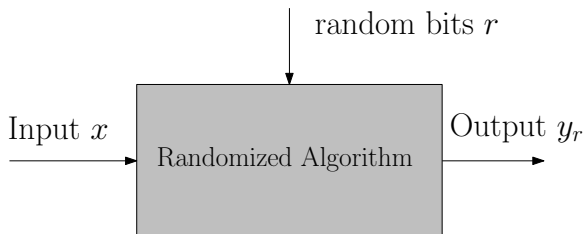
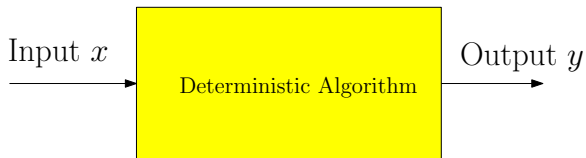
Part III

Randomization

Randomized Algorithms



Randomized Algorithms



Types of Randomized Algorithms

Typically one encounters the following types:

- 1 **Las Vegas randomized algorithms:** for a given input x output of *algorithm* is always correct but the *running time* is a *random variable*. In this case we are interested in analyzing the *expected* running time.

Types of Randomized Algorithms

Typically one encounters the following types:

- 1 **Las Vegas randomized algorithms:** for a given input x output of *algorithm* is *always correct* but the *running time* is a *random variable*. In this case we are interested in analyzing the *expected* running time.
- 2 **Monte Carlo randomized algorithms:** for a given input x the *running time* is *deterministic* but the *output* is *random*; correct with some probability. In this case we are interested in analyzing the *probability* of the correct output (and also the running time).
- 3 Algorithms whose running time and output may both be random.

Ping and find.

Consider a deterministic algorithm **A** that is trying to find an element in an array **X** of size **n**. At every step it is allowed to ask the value of one cell in the array, and the adversary is allowed after each such ping, to shuffle elements around in the array in any way it seems fit. For the best possible deterministic algorithm the number of rounds it has to play this game till it finds the required element is

- (A) $O(1)$
- (B) $O(n)$
- (C) $O(n \log n)$
- (D) $O(n^2)$
- (E) ∞ .

Ping and find randomized.

Consider an algorithm **randFind** that is trying to find an element in an array **X** of size **n**. At every step it asks the value of one random cell in the array, and the adversary is allowed after each such ping, to shuffle elements around in the array in any way it seems fit. This algorithm would stop in expectation after

- (A) $O(1)$
- (B) $O(\log n)$
- (C) $O(n)$
- (D) $O(n^2)$
- (E) ∞ .

steps.

Median

Consider the problem of finding an “approximate median” of an unsorted array $A[1..n]$: an element of A with rank between $n/4$ and $3n/4$.

- Finding an approximate median is not any easier than a proper median.
- $n/2$ elements of A qualify as approximate medians and hence a random element is good with probability $1/2$!

Part IV

Basics of Randomization

Discrete Probability Space

Definition

A discrete probability space is a pair (Ω, \Pr) consists of finite set Ω of **elementary events** and function $p : \Omega \rightarrow [0, 1]$ which assigns a probability $\Pr[\omega]$ for each $\omega \in \Omega$ such that $\sum_{\omega \in \Omega} \Pr[\omega] = 1$.

Example

An unbiased coin. $\Omega = \{H, T\}$ and $\Pr[H] = \Pr[T] = 1/2$.

Definition

Event is a collection of elementary events. The probability of an event $A \subset \Omega$, denoted by $\Pr[A]$, is $\sum_{\omega \in A} \Pr[\omega]$.

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Union Bound

For any two events \mathcal{E} and \mathcal{F} , we have that

$$\Pr[\mathcal{E} \cup \mathcal{F}] \leq \Pr[\mathcal{E}] + \Pr[\mathcal{F}].$$

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Independence

Events A and B are called independent if

$$\Pr[A \cap B] = \Pr[A] \Pr[B].$$

Random Variables

Definition

Given a probability space (Ω, \Pr) a (real-valued) random variable X over Ω is a function $X : \Omega \rightarrow \mathbb{R}$.

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Definition (Expectation: Average of \mathbf{X} as per \Pr)

Expectation of \mathbf{X} , $\mathbf{E}[\mathbf{X}]$, is defined as $\sum_{\omega \in \Omega} \Pr[\omega] \mathbf{X}(\omega)$.

If \mathbf{S} is the set of all values that \mathbf{X} takes, then expectation can also be written as $\sum_{x \in \mathbf{S}} x \Pr[\mathbf{X} = x]$.

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Linearity of Expectation

Given two random variables \mathbf{X}_1 and \mathbf{X}_2 ,
 $\mathbf{E}[\mathbf{X}_1 + \mathbf{X}_2] = \mathbf{E}[\mathbf{X}_1] + \mathbf{E}[\mathbf{X}_2]$.

Independence of Random Variables

Random variables **X** and **Y** are said to be independent if

$$\forall x, y, \quad \Pr[\mathbf{X} = x \wedge \mathbf{Y} = y] = \Pr[\mathbf{X} = x] \cdot \Pr[\mathbf{Y} = y]$$

Multiplication

If **X** and **Y** are independent then $\mathbf{E}[\mathbf{XY}] = \mathbf{E}[\mathbf{X}] \mathbf{E}[\mathbf{Y}]$.

Part V

Randomized Quick Sort

Randomized QuickSort

Randomized QuickSort

- ① Pick a pivot element *uniformly at random* from the array.
- ② Split array into 3 subarrays: those smaller than pivot, those larger than pivot, and the pivot itself.
- ③ Recursively sort the subarrays, and concatenate them.

Analysis via Recurrence

- 1 Given array \mathbf{A} of size n , let $Q(\mathbf{A})$ be number of comparisons of randomized **QuickSort** on \mathbf{A} .
- 2 Note that $Q(\mathbf{A})$ is a random variable.
- 3 Let $\mathbf{A}_{\text{left}}^i$ and $\mathbf{A}_{\text{right}}^i$ be the left and right arrays obtained if:

Let \mathbf{X}_i be indicator random variable, which is set to **1** if the pivot is of rank i in \mathbf{A} , else zero.

$$Q(\mathbf{A}) = n + \sum_{i=1}^n \mathbf{X}_i \cdot \left(Q(\mathbf{A}_{\text{left}}^i) + Q(\mathbf{A}_{\text{right}}^i) \right).$$

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$$Q(\mathbf{A}) = n + \sum_{i=1}^n X_i \cdot \left(Q(\mathbf{A}_{\text{left}}^i) + Q(\mathbf{A}_{\text{right}}^i) \right).$$

Since each element of \mathbf{A} has probability exactly of $1/n$ of being chosen:

$$E[X_i] = \Pr[\text{pivot is the element with rank } i] = 1/n.$$

Independence of Random Variables

Lemma

Random variables X_i is independent of random variables $Q(A_{left}^i)$ as well as $Q(A_{right}^i)$, i.e.

$$\begin{aligned} E[X_i \cdot Q(A_{left}^i)] &= E[X_i] E[Q(A_{left}^i)] \\ E[X_i \cdot Q(A_{right}^i)] &= E[X_i] E[Q(A_{right}^i)] \end{aligned}$$

Proof.

This is because the algorithm, while recursing on $Q(A_{left}^i)$ and $Q(A_{right}^i)$ uses new random coin tosses that are independent of the coin tosses used to decide the first pivot. Only the latter decides value of X_i . □

Analysis via Recurrence

Let $T(n) = \max_{A: |A|=n} E[Q(A)]$ be the worst-case expected running time of randomized **QuickSort** on arrays of size n .

We have, for any A :

$$Q(A) = n + \sum_{i=1}^n x_i \left(Q(A_{\text{left}}^i) + Q(A_{\text{right}}^i) \right)$$

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$$\Rightarrow E[Q(A)] \leq n + \sum_{i=1}^n \frac{1}{n} (T(i-1) + T(n-i)).$$

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We derived:

$$E[Q(A)] \leq n + \sum_{i=1}^n \frac{1}{n} (T(i-1) + T(n-i)).$$

Note that above holds for any A of size n . Therefore

$$\max_{A: |A|=n} E[Q(A)] = T(n) \leq n + \sum_{i=1}^n \frac{1}{n} (T(i-1) + T(n-i)).$$

Solving the Recurrence

$$T(n) \leq n + \sum_{i=1}^n \frac{1}{n} (T(i-1) + T(n-i))$$

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Lemma

$T(n) = O(n \log n)$.

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Lemma

$T(n) = O(n \log n)$.

Proof.

(Guess and) Verify by induction. □

Part VI

Inequalities

Markov's Inequality

Markov's inequality

Let \mathbf{X} be a **non-negative** random variable over a probability space (Ω, \mathbf{Pr}) . For any $\mathbf{a} > 0$,

$$\mathbf{Pr}[\mathbf{X} \geq \mathbf{a}] \leq \frac{\mathbf{E}[\mathbf{X}]}{\mathbf{a}}$$

Chebyshev's Inequality

Variance

Variance of \mathbf{X} is the measure of how much does it deviate from its mean value. Formally,

$$\mathbf{Var}(\mathbf{X}) = \mathbf{E}[(\mathbf{X} - \mathbf{E}[\mathbf{X}])^2] = \mathbf{E}[\mathbf{X}^2] - \mathbf{E}[\mathbf{X}]^2$$

Chebyshev's Inequality

Given $\mathbf{a} \geq 0$, $\Pr[|\mathbf{X} - \mathbf{E}[\mathbf{X}]| \geq \mathbf{a}] \leq \frac{\mathbf{Var}(\mathbf{X})}{\mathbf{a}^2}$

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Chebyshev's Inequality

Given $\mathbf{a} \geq 0$, $\mathbf{Pr[|X - E[X]| \geq a]} \leq \frac{\mathbf{Var(X)}}{\mathbf{a^2}}$

If \mathbf{X} and \mathbf{Y} are independent then $\mathbf{Var(X + Y) = Var(X) + Var(Y)}$.

Chebyshev's Inequality: Under Mutual Independence

Let $\mathbf{X}_1, \dots, \mathbf{X}_k$ be k independent random variables such that, for each $i \in [1, k]$, \mathbf{X}_i equals $\mathbf{1}$ with probability \mathbf{p}_i , and $\mathbf{0}$ with probability $(1 - \mathbf{p}_i)$. Let $\mathbf{X} = \sum_{i=1}^k \mathbf{X}_i$ and $\mu = \mathbf{E}[\mathbf{X}] = \sum_i \mathbf{p}_i$.

$$\mathbf{Var}(\mathbf{X}) \leq \mu \Rightarrow \mathbf{Pr}[|\mathbf{X} - \mu| \geq \mathbf{a}] \leq \frac{\mathbf{Var}(\mathbf{X})}{\mathbf{a}^2} < \frac{\mu}{\mathbf{a}^2}$$

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For $\delta > \mathbf{0}$, $\mathbf{Pr}[\mathbf{X} \geq (1 + \delta)\mu] \leq \frac{1}{\delta^2 \mu}$

For $\mathbf{0} < \delta < \mathbf{1}$, $\mathbf{Pr}[\mathbf{X} \leq (1 - \delta)\mu] \leq \frac{1}{\delta^2 \mu}$

Chernoff Bound

Let $\mathbf{X}_1, \dots, \mathbf{X}_k$ be k independent random variables such that, for each $i \in [1, k]$, \mathbf{X}_i equals $\mathbf{1}$ with probability \mathbf{p}_i , and $\mathbf{0}$ with probability $(\mathbf{1} - \mathbf{p}_i)$. Let $\mathbf{X} = \sum_{i=1}^k \mathbf{X}_i$ and $\mu = \mathbf{E}[\mathbf{X}] = \sum_i \mathbf{p}_i$. For any $\mathbf{0} < \delta < \mathbf{1}$, it holds that:

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$$\Pr[\mathbf{X} \geq (\mathbf{1} + \delta)\mu] \leq e^{-\frac{\delta^2 \mu}{3}} \text{ and } \Pr[\mathbf{X} \leq (\mathbf{1} - \delta)\mu] \leq e^{-\frac{\delta^2 \mu}{2}}$$

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Tighter bound

For any $\delta > 0$, $\Pr[\mathbf{X} \geq (\mathbf{1} + \delta)\mu] \leq \left(\frac{e^\delta}{(1+\delta)^{(1+\delta)}} \right)^\mu$

$$\Pr[\mathbf{X} \leq (\mathbf{1} - \delta)\mu] \leq \left(\frac{e^{-\delta}}{(1-\delta)^{(1-\delta)}} \right)^\mu$$

Problem: Approximate Median

Suppose you are presented with a very large set S of real numbers, and you would like to approximate the median of these numbers by sampling. You may assume all numbers in S are distinct. Let $|S| = n$. We say x is an ϵ -approximate median of S if at least $(1/2 - \epsilon)n$ are less than x and at least $(1/2 - \epsilon)n$ are greater than x . Consider an algorithm that samples $S' \subseteq S$ u.a.r. and outputs a median of S' . Show that there is an absolute constant c , independent of n , if the sample size is c , then with probability $1 - \delta$ the number returned will be ϵ -approximate median.

