EE360C: ALGORITHMS DYNAMIC PROGRAMMING

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Introduction

Weighted Interval Scheduling

Programming Generalities

Segmented Least Squares

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Greedy

Build up a solution incrementally, myopically optimizing some local criterion

Divide and Conquer

Break up a problem into two (or more) subproblems, solve each subproblem independently, and combine the solutions to the subproblems to form a solution to the original problem

Dynamic Programming

Break up a problem into a series of overlapping subproblems and build up solutions to larger and larger subproblems.

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Bellman pioneered the systematic study of dynamic programming in the 1950s.

Etymology

- Dynamic programming = planning over time
- Secretary of Defense at the time was hostile to mathematical research
- Bellman sought an impressive name to avoid confrontation
 - "something not even a Congressman could object to"

Dynamic Programming is an algorithm design technique that, like divide and conquer, relies on solutions to sub problems to solve a problem

- in dynamic programming, however, the subproblems are not independent
- a dynamic programming algorithm solves each of these subproblems just once, saving time in comparison to a traditional divide and conquer approach

Dynamic programming is commonly used for **optimization problems** in which many possible solutions exist; the algorithm helps us find one of the possibilities

- every solution to the problem has a value
- the algorithm helps us find a solution with the optimal value

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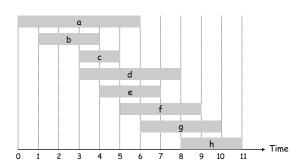
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Dynamic programming algorithms typically involve the following steps:

- characterize the structure of an optimal solution
- recursively define the value of an optimal solution
- compute the value of an optimal solution in a bottom-up fashion
- construct an optimal solution from the computed information

The weighted interval scheduling problem:

- Job j starts at s_j , finishes at f_j , and has weight or value v_i
- Two jobs are *compatible* if they do not overlap
- Goal: find maximum weight subset of mutually compatible jobs



UNWEIGHTED INTERVAL SCHEDULING

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Questions

Recall that greedy algorithm that works perfectly (optimally) if all the weights are 1. The greedy algorithm can fail spectacularly if arbitrary weights are allowed.

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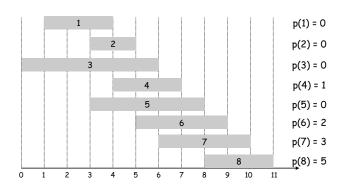
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Ouestions

Label (and sort) jobs by finishing time ($f_1 \le f_2 \le ... \le f_n$). Define p(j) to be the largest index i < j such that job i is compatible with job j.



OPT(j) is the value of the optimal solution to the problem consisting of job request 1, 2, ..., j.

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- Case 1: OPT selects job *j*.
 - then OPT can't include any of the jobs in the range p(j) + 1, p(j) + 2, ..., j 1
 - OPT must include the optimal solution to the problem consisting of remaining compatible jobs $1, 2, \dots, p(j)$
- Case 2: OPT does not select job *j*.
 - OPT must include the optimal solution to the problem consisting of remaining compatible jobs $1, 2, \ldots j-1$

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$$OPT(j) = \begin{cases} 0 & \text{if } j = 0\\ \max\{v_j + OPT(p(j)), OPT(j-1) & \text{otherwise} \end{cases}$$

```
Input: n, s_1,...,s_n, f_1,...,f_n, v_1,...,v_n

Sort jobs by finish times so that f_1 \le f_2 \le ... \le f_n.

Compute p(1), p(2), ..., p(n)

Compute-Opt(j) {
   if (j = 0)
     return 0
   else
     return max(v_j + Compute-Opt(p(j)), Compute-Opt(j-1))
}
```

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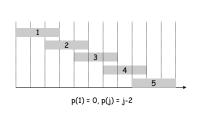
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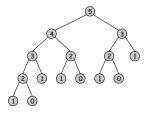
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Ouestions

The recursive algorithm fails spectacularly because of redundant sub-problems (it's an exponential algorithm)

 the number of recursive calls for a family of "layered" instances grows like the Fibonacci sequence





Memoization: store the results of each sub-problem in a cache; look them up as needed.

```
Input: n, s<sub>1</sub>,...,s<sub>n</sub>, f<sub>1</sub>,...,f<sub>n</sub>, v<sub>1</sub>,...,v<sub>n</sub>

Sort jobs by finish times so that f<sub>1</sub> ≤ f<sub>2</sub> ≤ ... ≤ f<sub>n</sub>.

Compute p(1), p(2), ..., p(n)

for j = 1 to n
    M[j] = empty ← global array

M[j] = 0

M-Compute-Opt(j) {
    if (M[j] is empty)
        M[j] = max(w<sub>j</sub> + M-Compute-Opt(p(j)), M-Compute-Opt(j-1))
    return M[j]
}
```

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Ouestions

What is the running time of the memoized algorithm? $O(n \lg n)$

- Sort by finish time: $O(n \lg n)$
- Computing $p(\cdot)$: O(n) after sorting by start time
- M-Compute-Opt(j): each invocation takes O(1) time and either:
 - returns an existing value M[j]
 - fills in one new entry M[j] and makes two recursive calls
- Define a progress measure Φ as the number of nonempty entries in M[].
 - initially $\Phi = 0$; throughout $\Phi \le n$
 - a call to M-Compute-Opt (j) increases Φ by at most 1; Φ cannot be increased more than n times; there are at most 2n recursive calls
- so the overall running time of M-Compute-Opt (n) is O(n)

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This algorithm has only computed the optimal value (i.e., the weight of the optimal schedule). What if we want the solution itself?

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```
Run M-Compute-Opt(n)
Run Find-Solution(n)

Find-Solution(j) {
   if (j = 0)
      output nothing
   else if (v<sub>j</sub> + M[p(j)] > M[j-1])
      print j
      Find-Solution(p(j))
   else
      Find-Solution(j-1)
}
```

To do this via dynamic programming, basically, we unwind the recursion...

```
Input: n, s_1, ..., s_n, f_1, ..., f_n, v_1, ..., v_n

Sort jobs by finish times so that f_1 \le f_2 \le ... \le f_n.

Compute p(1), p(2), ..., p(n)

Iterative-Compute-Opt {

M[0] = 0

for j = 1 to n

M[j] = \max(v_j + M[p(j)], M[j-1])
}
```

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WHEN DYNAMIC PROGRAMMING APPLIES

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Questions

For an optimization problem to be a good candidate for dynamic programming, it should exhibit:

- optimal substructure
- overlapping problems

A problem exhibits the optimal substructure property if the optimal solution contains within it optimal solutions to subproblems

 be careful, though, optimal substructure may also make a problem a good candidate for a greedy solution

We use the optimal substructure by then solving the subproblems bottom-up, combining the subproblem solutions as we move up.

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Onestions

If the recursive solution is going to solve the same subproblem multiple times, then we're wasting computation

- this is in contrast to problems that generate unique subproblems at every split, for which divide and conquer is a good fit
- one alternative for overlapping subproblems is *memoization*
 - it's like dynamic programming, but it fills the table in using a top-down approach that's more like traditional recursion
 - you create a table to use in a recursive algorithm, and the recursive calls fill in the table as they return
 - results can then be just computed once and looked up thereafter
- more traditional dynamic programming is "bottom up"

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Least Squares

- Foundational problem in statistic and numerical analysis
- Given n points in the plane $(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)$, find a line y = ax + b that minimizes the sum of the squared error:

$$SSE = \sum_{i=1}^{n} (y_i - ax_i - b)^2$$

Solution

The solution from calculus tells us that the minimum error is achieved when:

$$a = \frac{n \sum_{i} x_{i} y_{i} - (\sum_{i} x_{i})(\sum_{i} y_{i})}{n \sum_{i} x_{i}^{2} - (\sum_{i} x_{i})^{2}}$$

and

$$b = \frac{\sum_{i} y_i - a \sum_{i} x_i}{n}$$

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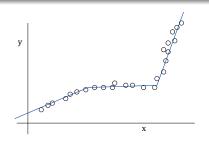
SEGMENTED LEAST SQUARES

Segmented Least Squares

- Points lie roughly on a sequence of several line segments
- Given n points in the plane $(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)$ with $x_1 < x_2 < \dots < x_n$, find a sequence of lines that minimizes f(x)

Question

What is a reasonable choice for f(x) to balance accuracy (goodness of fit) and parsimony (number of lines)



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The Tradeoff

We have to tradeoff the number of lines for the summed error

Find a sequence of lines that minimizes

- The sum of the sums of the squared errors *E* in each segment
- The number of lines *L*

This results in a tradeoff function E + cL for some constant c

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Notation

- $OPT(j) = \text{minimum cost for points } p_1, p_2, \dots, p_j$
- e(i,j) = minimum sum of squares for points p_i, p_{i+1}, \dots, p_j

To Compute OPT(j)

- Last segment uses points p_i, p_{i+1}, \dots, p_j for some i
- Cost = e(i, j) + c + OPT(i 1)

$$OPT(j) = \left\{ \begin{array}{ll} 0 & \text{if } j = 0 \\ \min_{1 \leq i \leq j} (e(i,j) + c + OPT(i-1)) & \textit{otherwise} \end{array} \right.$$

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```
INPUT: n, p<sub>1</sub>,...,p<sub>N</sub>, c

Segmented-Least-Squares() {
   M[0] = 0
   for j = 1 to n
      for i = 1 to j
      compute the least square error e<sub>ij</sub> for
      the segment p<sub>i</sub>,..., p<sub>j</sub>

for j = 1 to n
```

 $M[j] = min_{1,2,2,3} (e_{ij} + c + M[i-1])$

Running Time

return M[n]

• Bottleneck: computing e(i, j) for $O(n^2)$ pairs, O(n) per pair using the previous formula

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Knapsack Problem

- Given n objects and a "knapsack"
- Item *i* weighs $w_i > 0$ kilograms and has value $v_i > 0$
- Knapsack has a capacity of W kilograms
- Goal: fill the knapsack so as to maximize the total value

Greedy?

Remember the greedy algorithm we explored was not optimal...

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Ouestions

Definition: OPT(i) is the max profit for the subset of items $1,\ldots,i$

- Case 1: OPT does not select item i
 - *OPT* selects the best of $1, 2, \dots, i-1$
- Case 2: OPT selects item i
 - accepting item *i* does not immediately imply that we will have to reject other items
 - without knowing what other items were selected before i, we don't even know if we have enough room for i

Conclusion

We need more subproblems!

Definition: OPT(i, w) is the max profit of items $1, \dots, i$ with weight limit w

- Case 1: OPT does not select item i
 - *OPT* selects best of $\{1, 2, \dots, i-1\}$ using weight limit w
- Case 2: *OPT* selects item *i*
 - new weight limit $w w_i$
 - *OPT* selects best of $\{1, 2, ..., i 1\}$ using this new weight limit

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$$OPT(i, w) = \left\{ \begin{array}{ll} 0 & \text{if } i = 0 \\ OPT(i-1, w) & \text{if } w_i > w \\ \max(OPT(i-1, w), v_i + OPT(i-1, w-w_i)) & \text{otherwise} \end{array} \right.$$

```
Input: n, w_1, \dots, w_N, v_1, \dots, v_N
for w = 0 to W
   M[0, w] = 0
for i = 1 to n
   for w = 1 to W
       if (w_i > w)
          M[i, w] = M[i-1, w]
      else
          M[i, w] = \max \{M[i-1, w], v_i + M[i-1, w-w_i]\}
return M[n, W]
```

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Running Time

- $\theta(nW)$
- But I swear I've heard that the knapsack algorithm is NP-Complete!
- What does that mean?
- How is that possible? Didn't I just write down a polynomial time algorithm?
- Nope. It's "pseudo-polynomial"

An Approximation Algorithm

There is a polynomial time algorithm that produces a feasible solution whose value is within 0.01% of the optimal.

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Ouestions

You are given k denominations of coins, $d_1, d_2, \dots d_k$ (all integers). Assume $d_1 = 1$ so it is always possible to make change for any amount of money. We want to find an algorithm that makes change for an amount of money n using as few coins as possible.

First Question

The coin changing problem exhibits optimal substructure. Consider any optimal solution to make change for n cents using our k denominations of coins. Consider breaking that solution into two different pieces along any coin boundary. Suppose that one half of the solution amounts to b cents and the other half to n-b cents. Then the solution to the each half must be an optimal way to make change for b cents (or n-b cents) using the same k denominations of coins. **Prove this.**

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You are given k denominations of coins, $d_1, d_2, \ldots d_k$ (all integers). Assume $d_1 = 1$ so it is always possible to make change for any amount of money. We want to find an algorithm that makes change for an amount of money n using as few coins as possible.

Second Question

Let C[p] be the minimum number of coins of the k denominations that sum to p cents. **Recursively define the value of the optimal solution.** To get you started, there must exist some "first coin" d_i , where $d_i \leq p$.

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You are given k denominations of coins, $d_1, d_2, \ldots d_k$ (all integers). Assume $d_1 = 1$ so it is always possible to make change for any amount of money. We want to find an algorithm that makes change for an amount of money n using as few coins as possible.

Third Question

Provide an algorithm (e.g., pseudocode) to compute the value of the optimal solution bottom up. Provide another algorithm to construct the optimal solution from the computed information.

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You are given k denominations of coins, $d_1, d_2, \ldots d_k$ (all integers). Assume $d_1 = 1$ so it is always possible to make change for any amount of money. We want to find an algorithm that makes change for an amount of money n using as few coins as possible.

The Algorithm to Reconstruct the Solution

MAKECHANGE(S, d, n)

1 while n > 0

2 Print S[n]

 $n \leftarrow n - d[S[n]]$

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THE COIN CHANGING PROBLEM

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The Problem

You are given k denominations of coins, $d_1, d_2, \ldots d_k$ (all integers). Assume $d_1 = 1$ so it is always possible to make change for any amount of money. We want to find an algorithm that makes change for an amount of money n using as few coins as possible.

Fourth Question

What is the running time of your algorithm?

Running Time

CHANGE is $\Theta(nk)$ (which is pseudo-polynomial in k, the input size and dependent on n, just like knapsack was dependent on W). MAKECHANGE is O(n) since n is reduced by at least 1 in every iteration of the **while** loop. The total space requirement is $\Theta(n)$

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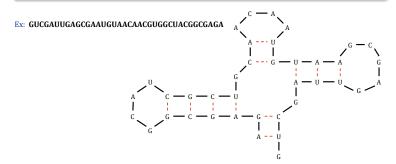
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Secondary Structure

RNA is single stranded, so it tends to loop back and form *base pairs* with itself. This structure is often essential to the function of the molecule. Legal base pairs are (A, U) or (C, G).



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Ouestions

- Watson and Crick: *S* is a matching and each pair in *S* is a Watson-Crick component (matching A to U or C to G)
- No Sharp Turns: The ends of each pair are separated by at least 4 intervening bases. That is, if $(b_i, b_j) \in S$, then i < j 4.
- Non-Crossing: If (b_i, b_j) and (b_k, b_l) are two pairs in S, the we cannot have i < k < j < l

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Ougetions

Free Energy

The usual hypothesis is that an RNA molecule will form the secondary structure that has the optimum total *free energy*, which we approximate by the number of base pairs.

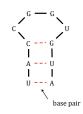
Goal

Given an RNA molecule $B = b_1b_2 \dots b_n$, find a secondary structure S that maximizes the number of base pairs.

RNA SECONDARY STRUCTURE: EXAMPLES

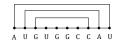
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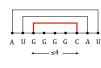
Examples.

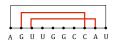












ok

sharp turn

crossing

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Notation

OPT(i, j) is the maximum number of base pairs in a secondary structure of the substring $b_i b_{i+1} \dots b_j$

- Case 1: if $i \ge j 4$
 - OPT(i, j) = 0 by the no sharp turns condition
- Case 2: Base b_i is not involved in a pair.
 - OPT(i, j) = OPT(i, j 1)
- Case 3: Base b_i pairs with b_t for some $i \le t < j 4$
 - The non crossing constraint decouples the resulting subproblems
 - $OPT(i.j) = 1 + \max_t (OPT(i, t 1) + OPT(t + 1, j 1))$ such that $i \le t < j 4$ and (b_j, b_t) is a Watson and Crick pair

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BOTTOM UP DYNAMIC PROGRAMMING OVER INTERVALS

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Question

What order to solve the subproblems?

Answer

Do the shortest intervals first.

```
RNA (b<sub>1</sub>,...,b<sub>n</sub>) {
   for k = 5, 6, ..., n-1
   for i = 1, 2, ..., n-k
        j = i + k
        Compute M[i, j]
   return M[1, n] using recurrence
}
```

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DYNAMIC PROGRAMMING SUMMARY

Recipe

- Characterize the structure of the problem
- Recursively define the value of an optimal solution
- Compute the value of the optimal solution
- Construct the optimal solution from computed information

Dynamic Programming Techniques

- Binary choice: weighted interval scheduling.
- Multi-way choice segmented least squares.
- Adding a new variable: knapsack
- Dynamic programming over intervals: RNA secondary structure

Top-down vs. Bottom-up?

Different people have different intuitions

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• useful in finding similar proteins, etc.

Dynamic Programming

Given a sequence $X = \langle x_1, x_2, \dots, x_m \rangle$, a sequence $Z = \langle z_1, z_2, \dots, z_k \rangle$ is a **subsequence** of X if there exists a strictly increasing sequence $\langle i_1, i_2, \dots, i_k \rangle$ of indices of X such that for all $j = 1, 2, \dots, k$, $x_{i_i} = z_j$.

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(For example $Z = \langle B, C, D, B \rangle$ is a subsequence of $X = \langle A, B, C, B, D, A, B \rangle$.)

Our goal, is, given two sequences *X* and *Y*, find the longest common subsequence of the two sequences.

CHARACTERIZING THE LONGEST COMMON SUBSEQUENCE

Theorem: Optimal Substructure of an LCS

Let $X = \langle x_1, x_2, \dots, x_m \rangle$ and $Y = \langle y_1, y_2, \dots, y_n \rangle$ be sequences and let $Z = \langle z_1, z_2, \dots, z_k \rangle$ be any LCS of X and Y.

- ① If $x_m = y_n$, then $z_k = x_m = y_n$ and Z_{k-1} is an LCS of X_{m-1} and Y_{n-1} .
- ② If $x_m \neq y_n$, then $z_k \neq x_m$ implies that Z is an LCS of X_{m-1} and Y.
- 3 if $x_m \neq y_n$, then $z_k \neq y_n$ implies that Z is an LCS of X and Y_{n-1}

Proof (Part I)

If $z_k \neq x_m$, then we could append $x_m = y_n$ to Z to get a common subsequence with length k+1, which would contradict the premise. So $z_k = x_m = y_n$. And therefore Z_{k-1} , the prefix of Z with z_k removed is a longest common subsequence of X_{m-1} and Y_{n-1} .

Proof (Parts II and III)

If $z_k \neq x_m$, then Z is a common subsequence of X_{m-1} and Y. If there were a common subsequence W of X_{m-1} and Y with length greater than k, then W would also be a subsequence of X and Y, which contradicts the premise that Z was an LCS.

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If $x_m = y_n$, there is one subproblem to solve: find the LCS of X_{m-1} and of Y_{n-1} . If $x_m \neq y_n$, then there are two subproblems to solve: find the LCS of X_{m-1} and Y and the LCS of X and Y_{n-1} .

$$c[i,j] = \begin{cases} 0 & \text{if } i = 0 \text{ or } j = 0 \\ c[i-1,j-1] + 1 & \text{if } i,j > 0 \text{ and } x_i = y_j \\ \max(c[i,j-1],c[i-1,j]) & \text{if } i,j > 0 \text{ and } x_i \neq y_j \end{cases}$$

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Ougotions

We use dynamic programming to compute solutions to the $\Theta(mn)$ subproblems in a bottom-up fashion

- we create a table c[0..m,0..n] and compute its entries in row major order (filling in the first row, left to right, then the second row, etc.)
- we keep a second table b[0..m,0..n] whose entries point to the table entry that corresponds to the best subproblem for the problem b[i,j] (to help in reconstructing the optimal solution)

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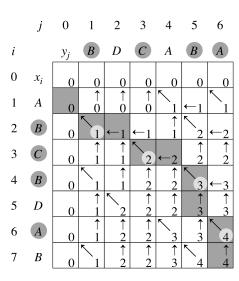
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```
LCS-LENGTH(X, Y)
       m \leftarrow \operatorname{length}[X]
   2 n \leftarrow \text{length}[Y]
   3 for i \leftarrow 1 to m
                do c[i, 0] \leftarrow 0
        for i \leftarrow 1 to n
   6
                do c[0,j] \leftarrow 0
        for i \leftarrow 1 to m
   8
                do for j \leftarrow 1 to n
                            do if x_i = y_i
  10
                                     then c[i, j] \leftarrow c[i - 1, j - 1] + 1
  11
                                             b[i,j] \leftarrow \nwarrow
  12
                                     else if c[i-1,j] \ge c[i,j-1]
  13
                                                then c[i, j] \leftarrow c[i-1], j
  14
                                                        b[i,j] \leftarrow \uparrow
  15
                                                else c[i,j] \leftarrow c[i,j-1]
  16
                                                        b[i,j] \leftarrow \leftarrow
        return c and b
```

AN EXAMPLE





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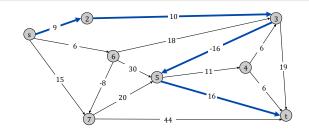
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Shortest Path Problem

Given a direct graph G = (V, E) with the edge weights c_{vw} (which can include negative edge weights), find the shortest path from node s to node t.

Example

Nodes represent agents in a financial setting and c_{vw} is the cost of a transaction in which we buy from agent v and sell immediately to w



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NEGATIVE COST CYCLES

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Observation

If some path from *s* to *t* contains a negative cost cycle, there does not exist a shortest *s-t* path; otherwise there exists a path that is a simple path.

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Definition: OPT(i, v) is the length of the shortest v-t path P using at most i edges.

- Case 1: P uses at most i 1 edges
 - OPT(i, v) = OPT(i 1, v)
- Case 2: P uses exactly i edges
 - if (v, w) is the first edge, the *OPT* uses (v, w), and then selects the best w-t path using at most i-1 edges

$$OPT(i,v) = \begin{cases} 0 & \text{if } i = 0 \\ \min(OPT(i-1,v), \min_{(v,w) \in E}(OPT(i-1,w) + c_{vw}))) & \text{otherwise} \end{cases}$$

IMPLEMENTATION

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Questions

See the book for the Bellman-Ford algorithm implementation and a discussion of the running time.

DISTANCE VECTOR ROUTING

Communication Network

- nodes = routers
- edges = direct communication link
- cost of edge = delay on link (which is naturally non-negative, but we use Bellman-Ford algorithm anyway!)

Dijkstra's Algorithm

Requires global information of network

Bellman-Ford

Uses only local knowledge of neighboring nodes

Synchronization

We don't expect the routers to run in lock-step; Bellman-Ford can tolerate asynchronous routing updates and can be shown to still converge on the correct shortest path.

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DISTANCE VECTOR PROTOCOL

Distance Vector Protocol

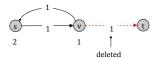
- Each router maintains a vector of shortest path lengths to every other node (distances) and the first hop on each path (directions)
- Algorithm: each router performs *n* separate computations, one for each potential destination node
- "Routing by rumor"

Examples

RIP, Xerox XNS RIP, Novell's IPX RIP, Cisco's IGRP, DEC's DNA Phase IV, AppleTalk's RTMP

Caveat

Edge costs may change during the algorithm; inconsistent router state can lead to the counting to infinity problem



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PATH VECTOR PROTOCOLS

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Link State Routing

- Each router stores the entire path (not just the distance and next hop)
- Based on Dijkstra's algorithm
- Avoids the counting to infinity problem
- Requires significantly more storage

Examples

Border Gateway Protocol (BGP), Open Shortest Path First (OSPF)

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