INF4130: Dynamic Programming September 6, 2016

- In the textbook: Ch. 9, and Section 20.5
 - We will use an example from Section 20.5 (Approximate string matching)
 - And we will study that example before we look at the theory
- The slides presented here have a different introduction to the theory than the textbook
 - This is done because the introduction in the textbook (built on its «principle of optimality») seems rather confusing.
 - Thus, some explanations for why the algorithms work will also be a little different than in the textbook.
- And the curriculum is the version used on these slides, not the introduction in the textbook.

Dynamic programming

Dynamic programming was formalised by Richard Bellmann (RAND Corporation) in the 1950'es.

- «programming» should here be understood as planning, or making decisions. It has nothing to do with writing code.
- "Dynamic" should indicate that it is a stepwise process.



Example: Given a string T and a shorter string P. Find strings «similar» to P in T (from Ch. 20.5)

We will first look at a similar problem, and leave the final step as an exercise This problem is similar to the one in 9.4: «Longest Common Subsequence»

A string *P* is a *k*-approximation of a string *T* if *T* can be converted to *P* by a sequence of maximum *k* of the following opertions:

Substitution One symbol in *T* is changed to another symbol.

Addition A new symbol is inserted somwhere in T.

Removing One symbol is removed from T.

The <u>Edit Distance</u>, ED(P,T), between two strings T and P is:

The smallest number of such operations needed to convert *T* to *P* (or *P* to *T*! Note that the definition is symmetric in *P* and *T*!)

Example.

```
logarithm → algarithm → algarithm → algorithm (Steps: +a, -o, a->o)

T

P
```

Thus ED("logarithm", "algorithm") = 3 (as there are no shorter way!)

An idea for finding the Edit Distance

To solve this problem we will use an interger matrix D[0:m, 0:n] as shown below, where we imagine that the string P = P[1:m] is placed downwords along the left side of D, and T = T[1:n] is placed above D from left to right (at corresponding indices).

(This is a slightly different use of indices than in Sectons 20.5 and 9.4).

Our plan is then to systematically fill in this table so that

D[i, j] = Edit Distance between the strings P[1:i] and T[1:j]

We will do this from «smaller» to «larger» index pairs (i, j), by taking column after column from left to right (but row after row from top to bottom would also work). The value we are looking for, ED(P, T), will then occur in D[m, n] when all entries are filled in.

The matrix *D*:

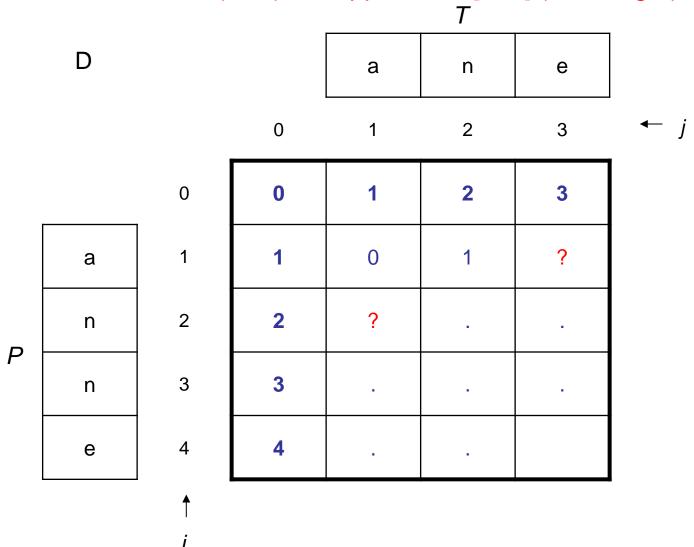
j -1

n

- Some obvious values for D[i, j] are given below
- How to find more values?
- The answer ED(P, T) will appear in D[4, 3] (lower right)

		•	, ,		T		
	D			а	n	е	
			0	1	2	3	← j
		0	0	1	2	3	
	а	1	1	?			
_	n	2	2				
Р	n	3	3				
	e	4	4				
		†					
		j					

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	D			а	n	е	
			0	1	2	3	← j
		0	0	1	2	3	
	а	1	1	0	1	2	
	n	2	2	?	?	?	
Р	n	3	3				
	e	4	4				
		<u></u>					
		i					

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		`	,		T	_ `	
	D			а	n	е	
			0	1	2	3	← j
		0	0	1	2	3	
	а	1	1	0	1	2	
	n	2	2	1	0	1	
P	n	3	3				
	e	4	4				
		†					
		i					

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		`	,		T	_ `	
	D			а	n	е	
			0	1	2	3	← j
		0	0	1	2	3	
	а	1	1	0	1	2	
_	n	2	2	1	0	1	
Р	n	3	3	2	1	1	
	e	4	4	3	2	1	
		† ;					•

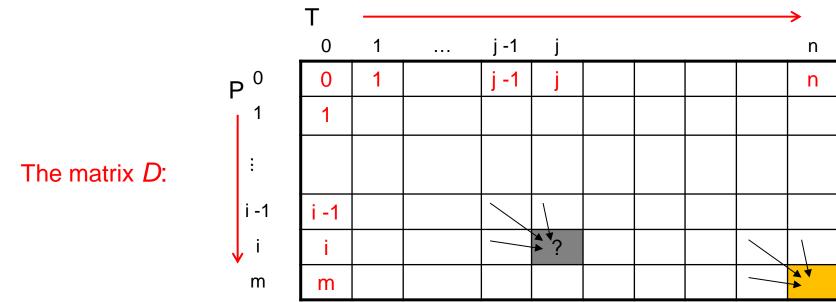
The «recurrence relation»

To be able to fill in this matrix we shall use that we have the relation indicated below between neighbouring entries in D. For proof see next slide.

Note that the value of D[i,j] only depends on entries in D with «smaller» index-pairs: D[i-1,j-1], D[i-1,j], and D[i,j-1]

$$D[i,j] = \begin{cases} D[i-1,j-1] & \text{hvis } P[i] = T[j] \\ \min \{ \underbrace{D[i-1,j-1]+1}_{\text{substitusj on}}, \underbrace{D[i-1,j]+1}_{\text{sletting } iP}, \underbrace{D[i,j-1]+1}_{\text{sletting } iP} \} & \text{ellers} \\ D[0,0] = 0, \quad D[i,0] = D[0,i] = i. \end{cases}$$

The equalities at the last line can be used to initialize the matrix (shown in red).



Proof of the recurrence relation

As indicated: The previous slides do not give a full proof for the equality in the recurrence relation (but only a «less than or equal»):

$$D[i,j] = \begin{cases} D[i-1,j-1] & \text{if } P[i] = T[j] \\ \min\{ D[i-1,j-1]+1, D[i-1,j]+1, D[i,j-1]+1 \} & \text{otherwise} \end{cases}$$
substitution addition to T Deletion from T

- This is because we can easily give D[i, j] steps to change P to T
- To prove the full equality will be an exercise this week.
- As a base for this proof we will assume that a minimal sequence of edit operations can always be written in the following form:

Here '+' = insertion, '-' = deletion, and '*' = substitution.

Figure for computing the edit distance (see program next slide)

We have the following expression for D[i, j]:

$$D[i,j] = \begin{cases} D[i-1,j-1] & \text{if } P[i] = T[j] \\ \min\{ \underbrace{D[i-1,j-1]+1}_{\text{substitution}}, \underbrace{D[i-1,j]+1}_{\text{addition to T}}, \underbrace{D[i,j-1]+1}_{\text{Deletion from T}} \} \text{ otherwise} \end{cases}$$

$$D[0,0] = 0$$
, $D[i,0] = D[0,i] = i$.

	0	1	 j -1	j	_		n
0	0	1	j -1	j			
1	1						
i							
i -1	i -1		/				
i	·i			*			\
m							*

When fully filled in, we will find the edit distance between T and P in D[m,n]

A program for computing the edit distance

```
function EditDistance ( P [1:m ], T [1:n ] )
     for i \leftarrow 0 to m do D[i, 0] \leftarrow i endfor
     for j \leftarrow 1 to n do D[0, j] \leftarrow j endfor
     for i \leftarrow 1 to m do
           for j \leftarrow 1 to n do
                  If P[i] = T[j] then
                       D[i,j] \leftarrow D[i-1,j-1]
                  else
                       D[i, j] \leftarrow \min \{ D[i-1, j-1] + 1, D[i-1, j] + 1, D[i, j-1] + 1 \}
                  endif
           endfor
                              Note that, after the initialization, we look at the pairs (i, j)
     endfor
                              in the following order (line after line at previous slide):
     return( D[ m, n ] )
```

end EditDistance

Note that, after the initialization, we look at the pairs (i, j) in the following order (line after line at previous slide):

(1,1) (1,2) ... (1,n)

(2,1) (2,2) ... (2,n)

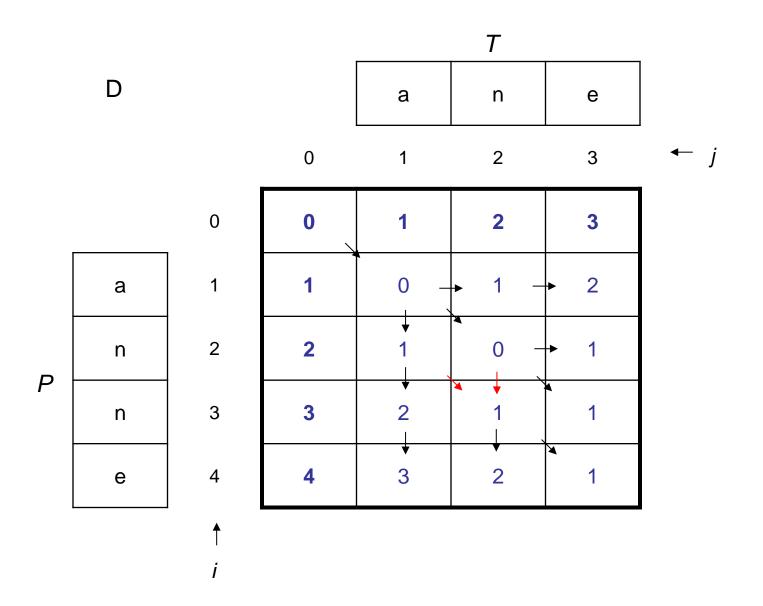
...

(m,1) ... (m,n)

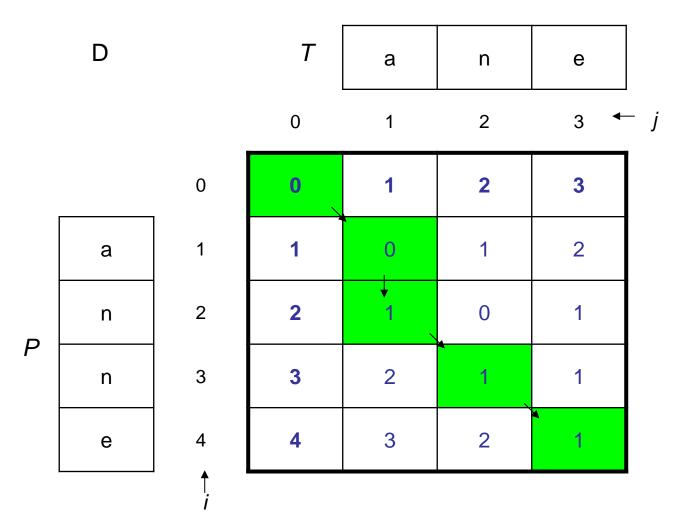
This is OK as this order ensures that the smaller instances are solved before they are needed to solve a larger instance. That is:

```
D[i-1, j-1], D[i-1, j] and D[i, j-1] are computed before D[i, j],
```

Our old example: Computing Edit distance



Finding the edit steps



Method: Follow the «path» used from the final entry back to [0,0]:

Diagonally, and P[i] = T[j]:

- No edit needed

Diagonally, and $P[i] \neq T[j]$:

- Substitution (Appear for D[3, 3])

Downwards:

- A letter is deleted from P

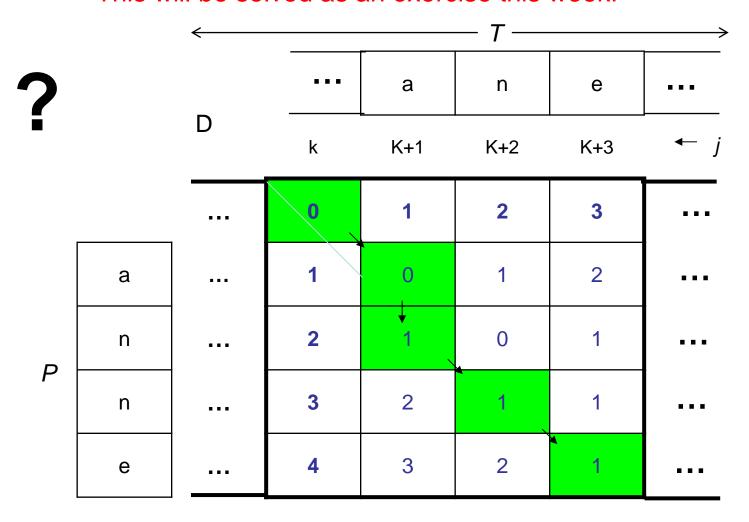
Towards the right:

- A letter is added to P (Appear for D[1, 2])

Until now we have computed the edit distance between two strings P and T

But what about searching for substrings U in T so that ED(P,U) is small, e.g, smaller than a certain given value?

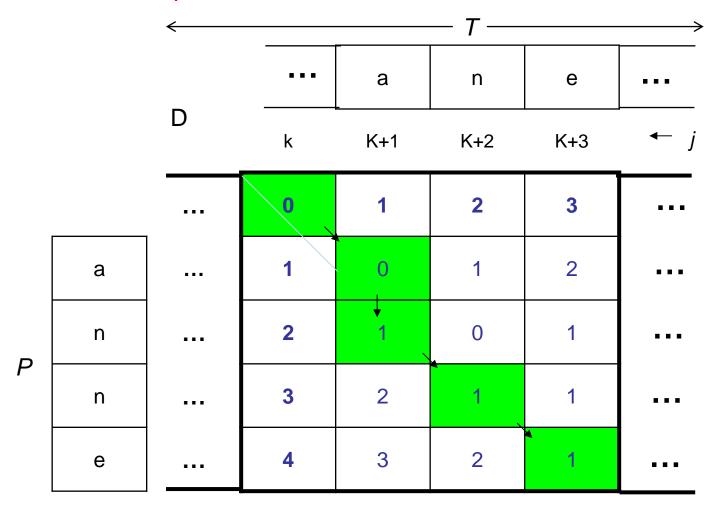
This will be solved as an exercise this week!



Relevance for research in genetics

Then T may be the full «genome» of one organism, and P a part of the genome of another.

Question: Does a sequence similar to P occur in T?

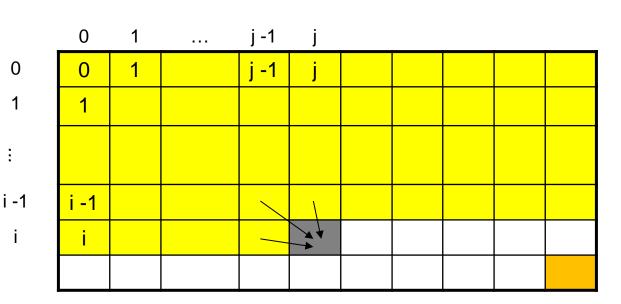


About Dynamic Programming in general

- Dynamic programming is typically used to solve optimazation problems P.
- The instances of the problem must be ordered from smaller to larger ones, and the smallest (or simplest) instances can usually easily be solved.
- For each problem instance I there is a set of instances I₁, I₂, ..., I_k, all smaller than I, so that we can find an (optimal) solution to I if we know the (optimal) solution of the I_i-problems.

In our example:

'The values of the yellow area is computed when the white value is to be computed



A general formulation of Dynamic Programming

Assume we have:

- A problem P with (usually infinitely) many instances I_1 , I_2 , $I_{3,...}$
- An integer $\langle size \rangle$ for each instance, where the $\langle simplest \rangle$ instances have a small size, usually 0 or 1. (In our example the size can be chosen as m+n)
- The solution to instance *I* is written *s*(*I*)
- For each I there is a set of instances J_1, \ldots, J_k called the *base of I*, written $B(I) = \{J_1, J_2, \ldots, J_k\}$ (where k may vary with I), and every J_k is smaller than I.
- We have a process/function *Combine* that takes as input an instance I, and the solutions s(J) to all J in B(I),

```
s(I) = Combine(I, s(J_1), s(J_2), ..., s(J_k))
```

This is called the «recurrence relation» of the problem.

- For an instance I, we can set up a sequence of instances $< L_0, L_1, ..., L_m = I >$ so that for all $p \le m$, all instances in $B(L_p)$ occur in the sequence before L_p .
- The sequence can be stored in a suitable table of reasonable size compared to the size of the instance *I*.

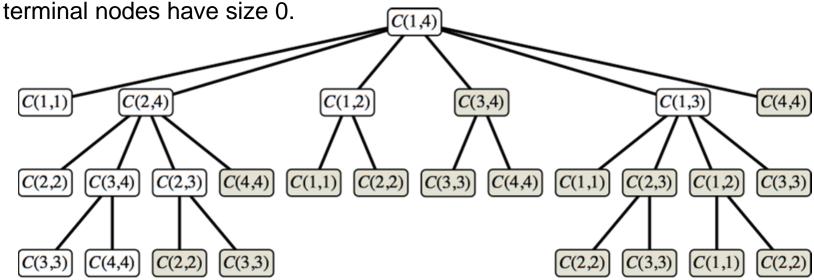
When to use dynamic programming?

- Dynamic programming is useful if the total number of smaller instances needed to solve an instance / (found recursively by the Bfunction) is so small that the answer to all of them can be stored in a suitable table.
- The main trick is to store the solutions in the table for later use. The real gain comes when each «smaller» entry is used a number of times.

	0	1	 j -1	j			
0	0	1	j -1	j			
1	1						
:							
i -1	i -1		//	\			
i	i		+	×			
					·		

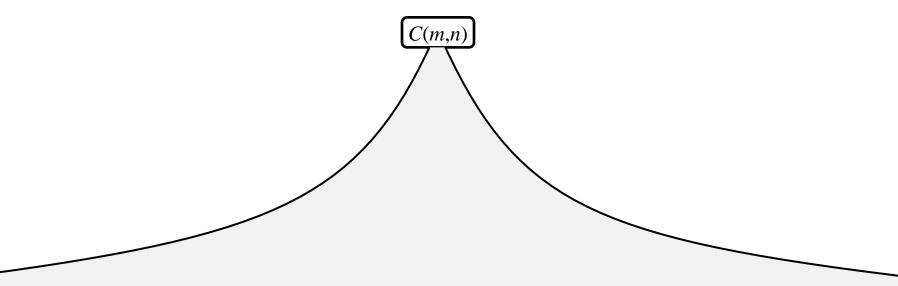
Another (slightly abstract) example

- As indicated on the previous slide, Dynamic Programming is useful if the solution to a certain instance is used in the solution of many instances of larger size.
- In the problem C below, an instance is given by some data (e.g two strings) and by two intergers *i* and *j*. Assume the corresponding instances are written C(i, j). Thus the solutions to the instances can be stored in a two-dimentional table with dimensions *i* and *j*.
- Assume that the size of an instanstance C(i, j) is j − i
- Below, the children of a node *N* indicate the instances of which we need the solution for computing the solution to *N* (which all have smaller size)
- Note that the solution to many instances, e.g. C(3,4), is used multiple times.
 This is required to make DP a preferable alternative! Also note that all terminal nodes have size 0.



Number of smaller solutions needed

- If the number of solutions to smaller instances needed to find the soluton to a certain instance is very big (e.g. exponential in the size of the instance), then the resulting DP algoritm will usually not be practical, as the table must be very big, and there will probably be little reuse of smaller solutions
- For Dynamic Programming to work well, this number should at most be polynomial in the size of the instance, and and usually rather small.



Sketch indicating a situation when the solution of one instance needs the solution of a large number of *different* smaller instances.

Two variants of dynamic programming: Bottom up (traditional) and top down (memoization)

Traditional Dynamic Programming (bottom up)

- Is traditionally performed bottom-up. All relevant smaller instances are solved first, and their solutions are stored in the table.
- This usually leads to very simple and often rapid programs.

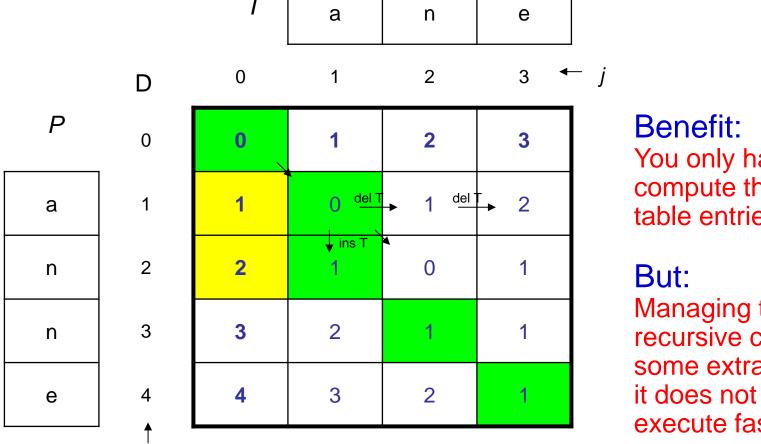
«Top-Down» Dynamic Programming

A drawback with (traditional) dynamic programming is that one usually solves a number of smaller instances that turn out not to be needeed for the actual (larger) instance that we are really interested in.

- We can instead start at the (large) instance we want to solve, and do the computation recursively top-down. Also here we put computed solutions in the table as soon as they are computed.
- Each time we need to solve an instance we first check whether it is already solved, and if so we only use the stored solution, otherwise we do recursive calls.
- The table entries then need a special marker «not computed», which also should be the initial value of the entries.

«Top-Down» dynamic programming: "Memoization" Method:

- 1. Start at the instance you want to solve, and ask recursively for the solution to the instnces needed
- 2. As soon as you have an answer, fill it into the table, and take it from there when the answer to the same instance is later needed.



You only have to compute the needed table entries

Managing the recursive calls take some extra time, so it does not always execute fastest.

Another example: Optimal Matrix Multiplication

Given the sequence $M_0, M_1, ..., M_{n-1}$ of matrices. We want to compute the product: $M_0 \cdot M_1 \cdot ... \cdot M_{n-1}$.

Note that for this multiplication to be meaningful the length of the rows in M_i must be equal to the length of the columns M_{i+1} for i = 0, 1, ..., n-2

Matrix multiplication is associative: $(A \cdot B) \cdot C = A \cdot (B \cdot C)$ But it is not symmetric, since $A \cdot B$ generally is different from $B \cdot A$

Thus, one can do the multiplications in different orders. E.g., with four matrices it can be done in the following five ways:

```
(M_0 \cdot (M_1 \cdot (M_2 \cdot M_3)))
(M_0 \cdot ((M_1 \cdot M_2) \cdot M_3))
((M_0 \cdot M_1) \cdot (M_2 \cdot M_3))
((M_0 \cdot (M_1 \cdot M_2)) \cdot M_3)
(((M_0 \cdot M_1) \cdot M_2) \cdot M_3)
```

The cost (the number of simple (scalar) multiplications) for these will usually vary a lot for the differnt alternatives. We want to find the one with as few scalar multiplications as possible.

Optimal matrix multiplication 2

Given two matrices A and B with dimentions:

```
A is a p \times q matrix,
B is a q \times r matrix.
```

The cost of computing $A \cdot B$ is $p \cdot q \cdot r$, and the result is a $p \times r$ matrix

Example

Compute $A \cdot B \cdot C$, where

A is a 10 \times 100 matrix, B is a 100 \times 5 matrix, and C is a 5 \times 50 matrix.

Computing $D = (A \cdot B)$ costs 5,000 and gives a 10 × 5 matrix. Computing $D \cdot C$ costs 2,500.

Total cost for $(A \cdot B) \cdot C$ is thus 7,500.

Computing $E = (B \cdot C)$ costs 25,000 and gives a 100 × 50 matrix. Computing $A \cdot E$ costs 50,000. Total cost for $A \cdot (B \cdot C)$ is thus 75,000.

We would indeed prefer to do it the first way!

Optimal matrix multiplication 3

Given a sequence of matrices M_0 , M_1 , ..., M_{n-1} . We want to find the cheapest way to do this multiplication (that is, an optimal paranthesization).

From the outermost level, the first step in a parenthesizaton is a partition into two parts: $(M_0 \cdot M_1 \cdot \ldots \cdot M_k) \cdot (M_{k+1} \cdot M_{k+2} \cdot \ldots \cdot M_{n-1})$ If we know the best parenthesizaton of the two parts, we can sum their cost and get the cost of the best parenthesizaton given that we have to use this *outermost* partition.

Thus, to find the best parenthesizaton of M_0 , M_1 , ..., M_{n-1} , we can simply look at all the n-1 possible outermost partitions (k = 0, 1, n-2), and choose the best. But we will then need the cost of the optimal parenthesizaton of a lot of instances of smaller sizes.

And we shall say that the *size* of the instance M_i , M_{i+1} , ..., M_j is j - i.

We therefore generally have to look at the best parenthesizaton of all intervals M_i , M_{i+1} , ..., M_j , in the order of growing sizes.

We will refer to the lowest possible cost for M_i , M_{i+1} , ..., M_j as $m_{i,j}$.

Optimal matrix multiplication 4

Let $d_0, d_1, ..., d_n$ be the dimensiones of the matrices $M_0, M_1, ..., M_{n-1}$, so that matrix M_i has dimension $d_i \times d_{i+1}$

As on the previous slide:

Let $m_{i,j}$ be the cost of an optimal parenthesizaton of M_i , M_{i+1} , ..., M_j . Thus the value we are interested in is $m_{0,n-1}$

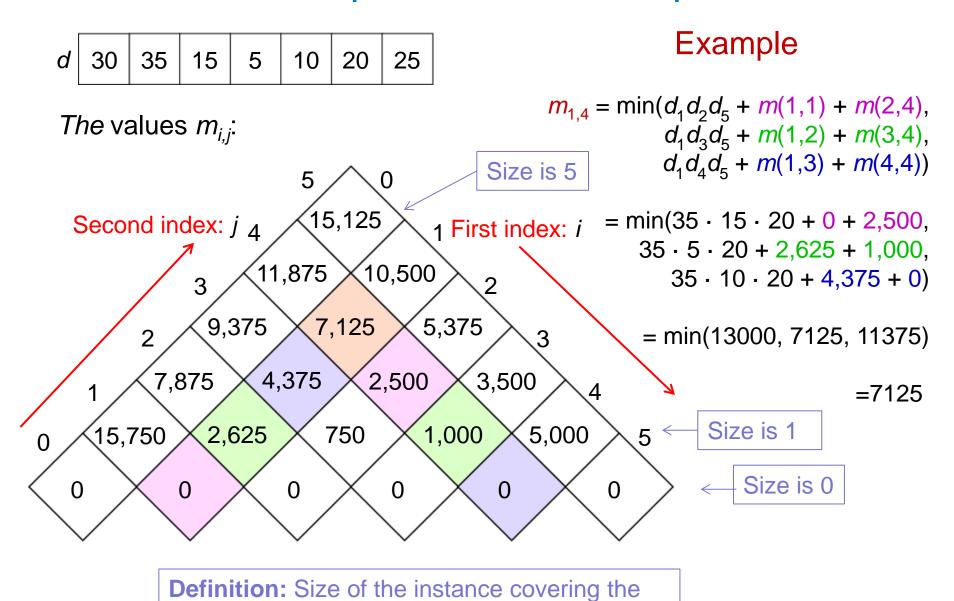
The recurrence relation for $m_{i,j}$ will be:

$$m_{i,j} = \min_{i \le k < j} \{ m_{i,k} + m_{k+1,j} + d_i d_{k+1} d_{j+1} \}, \text{ when } 0 \le i < j \le n-1$$

$$m_{i,i} = 0 \text{ when } 0 \le i \le n-1$$

Note that the values $m_{k,l}$ that we need for computing $m_{i,j}$ are all for smaller instances. That is: l - k < j - i.

The table: Optimal matrix multiplication



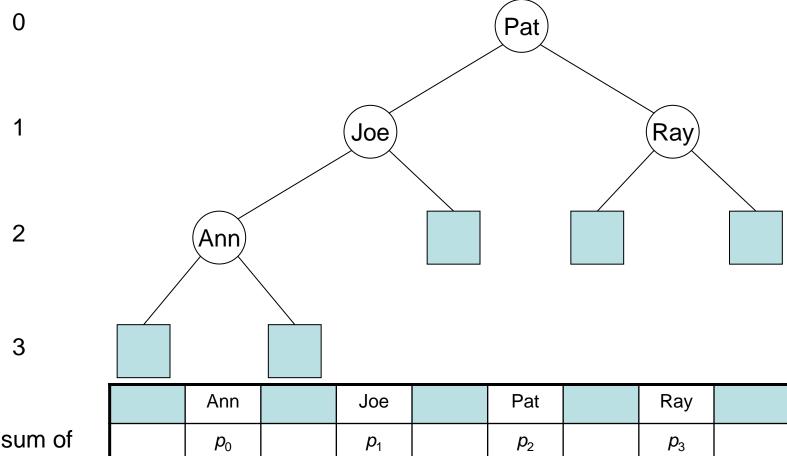
interval from pos. i to pos. j is i - i

Program: Optimal matrix multiplication

```
function OptimalParens( d[0 : n – 1] )
    for i \leftarrow 0 to n-1 do
            m[i, i] \leftarrow 0
    for diag \leftarrow 1 to n – 1 do
           for i \leftarrow 0 to n - 1 - diag do
                       i ← i + diag
                        m[i, j] \leftarrow \infty // Relative to the scalar values that can occur
                        for k \leftarrow i to j - 1 do
                                    q \leftarrow m[i, k] + m[k + 1, i] + d[i] \cdot d[k + 1] \cdot d[i + 1]
                                    if q < m[i, j] then
                                                m[i, i] \leftarrow q
                                               c[i,j] \leftarrow k
                                    endif
    return m[0, n – 1]
end OptimalParens
```

Yet another example: Optimal search trees

(Not in the curriculum for 2016)



The sum of the *p*'s and *q*'s is 1

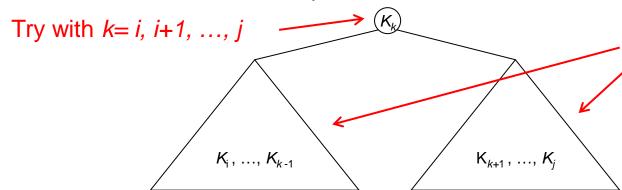
		Ann		Joe		Pat		Ray	
		p_0		p_1		p_2		p_3	
Ī	q_0		q_1		q_2		q_3		q_4
	3	3	3	2	2	1	2	2	2

Average search time: $3p_0 + 2p_1 + 1p_2 + 2p_3 + 3q_0 + 3q_1 + 2q_2 + 2q_3 + 2q_4$

Optimal search trees

(Not in the curriculum for 2016)

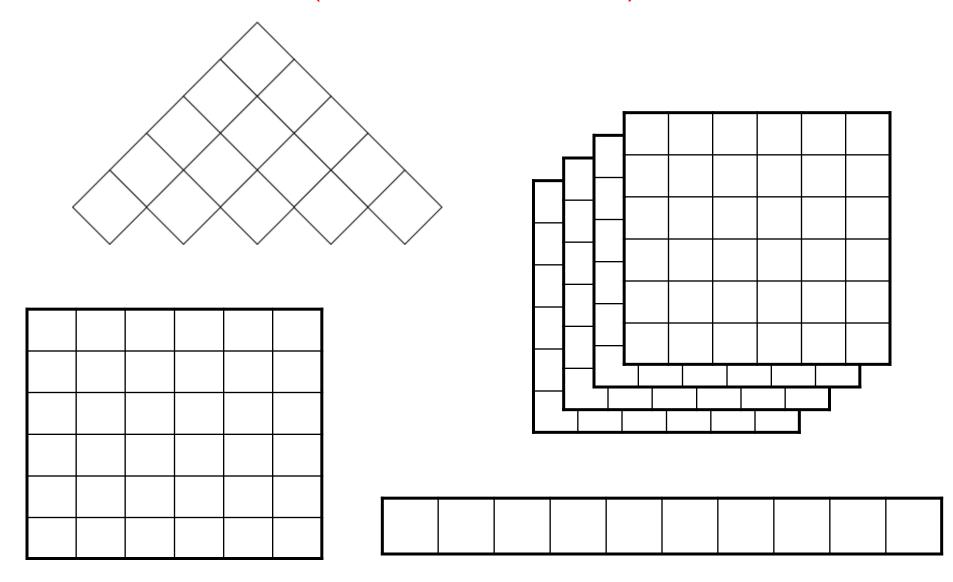
- To get a managable problem that still catches the essence of the general problem, we shall assume that all q-es are zero (that is, we never search for values not in the tree)
- A key to a solution is that a subtree in a search tree will always represent an interval of the values in the tree in sorted order (and that such an interval can be seen as an optimal seach instance in itself)
- Thus, we can use the same type of table as in the matrix multiplication case, where the value of the optimal tree over the values from intex i to index j is stored in A[i, j], and the size of such an instance is j i
- Then, for finding the optimal tree for an interval with values K_i , ..., K_j we can simply try with each of the values K_i , ..., K_j as root, and use the best subtrees in each of these cases (whose optimal values are already found).
- To compute the cost of the subtrees is slightly more complicated than in the matrix case, but is no problem.



The optimal values and form for these subtrees are already computed, when we here try with different values K_k at the root

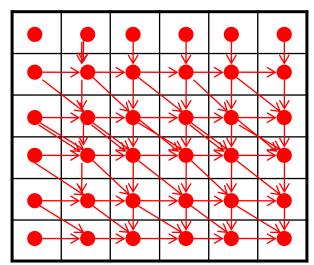
Dynamic programming in general:

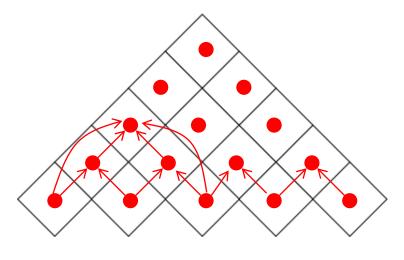
We fill in differnt types of tables «bottom up» (smallet instances first)



Dynamic programming Filling in the tables

- It is always safe to solve all the smaller instances before any larger ones, using the defined size of the instances.
- However, if we know what smaller instances are needed to solve a larger instance, we can deviate from the above. The important thing is that the smaller instances needed to solve a certain instance J is computed before we solve J.
- Thus, if we know the "dependency graph" of the problem (which must be cycle-free, see examples below), the important thing is to look at the instances in an order that conforms with this dependency. This freedom is often utilized to get a simple computation (see next slide).





Dynamic Programming using memoization

«Top-Down» dynamic programming (Memoization)

- A drawback with bottom up dynamic programming is that you might solve a lot of smaller instances whose answers are never used.
- We can instead do the computation recursively from the top, and store the (really needed) answers of the smaller instances in the same table as before. Then we can later find the answers in this table if we need the answer to the same instance once more.
- The reason we do not always use this technique is that recursion in itself can take a lot of time, so that a simple bottom up may be faster.
- For the recursive method to work, we need a flag «NotYetComputed» in each entry, and if this flag is set when we need that value, we compute it, and save the result (and turn off the flag, so the recursion from here will only be done once).
- The «NotYetComputed» flag must be set in all entries at the start of the algorithm.

Dynamic programming using memoization

- It is always safe to solve all the smaller instances before any larger ones, using the defined size of the instances.
- However, if we know what smaller instances are needed to solve a larger instance, we can deviate from that. The important thing is that the smaller instances needed to solve a certain instance J is computed before we start solving J.
- Thus, if we know the "dependency graph" of the problem (which must be cycle-free, see examples below), the important thing is to look at the instances in an order that conforms with this dependency. This freedom is often utilized to get a simple computation.

At most the entries with a green dot will have to be computed

