Dynamic Programming

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Topic Overview

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Overview of Serial Dynamic Programming

- Dynamic programming (DP) is used to solve a wide variety of discrete optimization problems such as scheduling, string-editing, packaging, and inventory management.
- Break problems into subproblems and combine their solutions into solutions to larger problems.
- In contrast to divide-and-conquer, there may be relationships across subproblems.

Dynamic Programming: Example

- Consider the problem of finding a shortest path between a pair of vertices in an acyclic graph.
- An edge connecting node *i* to node *j* has cost *c(i,j)*.
- The graph contains n nodes numbered 0,1,..., n-1, and has an edge from node i to node j only if i < j. Node 0 is source and node n-1 is the destination.
- Let f(x) be the cost of the shortest path from node 0 to node x.

$$f(x) = \left\{ egin{array}{cc} 0 & x = 0 \ \min_{0 \leq j < x} \{f(j) + c(j, x)\} & 1 \leq x \leq n-1 \end{array}
ight.$$

Dynamic Programming: Example



 A graph for which the shortest path between nodes 0 and 4 is to be computed.

$$f(4) = \min\{f(3) + c(3,4), f(2) + c(2,4)\}.$$

Dynamic Programming

- The solution to a DP problem is typically expressed as a minimum (or maximum) of possible alternate solutions.
- If *r* represents the cost of a solution composed of subproblems x₁, x₂,..., x_i, then *r* can be written as

$$r = g(f(x_1), f(x_2), \ldots, f(x_l)).$$

Here, *g* is the *composition function*.

• If the optimal solution to each problem is determined by composing optimal solutions to the subproblems and selecting the minimum (or maximum), the formulation is said to be a DP formulation.

Dynamic Programming: Example



The computation and composition of subproblem solutions to solve problem $f(x_8)$.

Dynamic Programming

- The recursive DP equation is also called the *functional* equation or optimization equation.
- In the equation for the shortest path problem the composition function is f(j) + c(j,x). This contains a single recursive term (f(j)). Such a formulation is called monadic.
- If the RHS has multiple recursive terms, the DP formulation is called polyadic.

Dynamic Programming

- The dependencies between subproblems can be expressed as a graph.
- If the graph can be levelized (i.e., solutions to problems at a level depend only on solutions to problems at the previous level), the formulation is called serial, else it is called non-serial.
- Based on these two criteria, we can classify DP formulations into four categories - serial-monadic, serialpolyadic, non-serial-monadic, non-serial-polyadic.
- This classification is useful since it identifies concurrency and dependencies that guide parallel formulations.

Serial Monadic DP Formulations

- It is difficult to derive canonical parallel formulations for the entire class of formulations.
- For this reason, we select two representative examples, the shortest-path problem for a multistage graph and the 0/1 knapsack problem.
- We derive parallel formulations for these problems and identify common principles guiding design within the class.

- Special class of shortest path problem where the graph is a weighted multistage graph of r + 1 levels.
- Each level is assumed to have *n* levels and every node at level *i* is connected to every node at level *i* + 1.
- Levels zero and *r* contain only one node, the source and destination nodes, respectively.
- The objective of this problem is to find the shortest path from S to R.



An example of a serial monadic DP formulation for finding the shortest path in a graph whose nodes can be organized into levels.

- The *ith* node at level *l* in the graph is labeled *v_i^l* and the cost of an edge connecting *v_i^l* to node *v_i^{l+1}* is labeled *c_{i,i}^l*.
- The cost of reaching the goal node R from any node vⁱ_i is represented by Cⁱ_i.
- If there are *n* nodes at level *I*, the vector $[C_0^{\ I}, C_1^{\ I, \dots, } C_{n-1}^{\ I}]^T$ is referred to as C^I . Note that $C_0 = [C_0^{\ 0}].$
- We have $C'_{i} = min \{ (c'_{i,j} + C'_{j+1}) \mid j \text{ is a node at level } l + 1 \}$

- Since all nodes v_j^{r-1} have only one edge connecting them to the goal node R at level r, the cost C_j^{r-1} is equal to $c_{j,r}^{r-1}$.
- We have:

$$\mathcal{C}^{r-1} = [c_{0,R}^{r-1}, c_{1,R}^{r-1}, \dots, c_{n-1,R}^{r-1}].$$

Notice that this problem is serial and monadic.

 The cost of reaching the goal node R from any node at level l is (0 < l < r - 1) is

$$\begin{split} C_0^l &= \min\{(c_{0,0}^l + C_0^{l+1}), (c_{0,1}^l + C_1^{l+1}), \dots, (c_{0,n-1}^l + C_{n-1}^{l+1})\},\\ C_1^l &= \min\{(c_{1,0}^l + C_0^{l+1}), (c_{1,1}^l + C_1^{l+1}), \dots, (c_{1,n-1}^l + C_{n-1}^{l+1})\},\\ &: \end{split}$$

 $C_{n-1}^{l} = \min\{(c_{n-1,0}^{l} + C_{0}^{l+1}), (c_{n-1,1}^{l} + C_{1}^{l+1}), \dots, (c_{n-1,n-1}^{l} + C_{n-1}^{l+1})\}.$

- We can express the solution to the problem as a modified sequence of matrix-vector products.
- Replacing the addition operation by minimization and the multiplication operation by addition, the preceding set of equations becomes:

$$\mathcal{C}^l = M_{l,l+1} \times \mathcal{C}^{l+1},$$

where C' and C'^{+1} are $n \ge 1$ vectors representing the cost of reaching the goal node from each node at levels *I* and I + 1.

 Matrix M_{I,I+1} is an n x n matrix in which entry (i, j) stores the cost of the edge connecting node i at level I to node j at level I + 1.

$$M_{l,l+1} = \begin{bmatrix} c_{0,0}^l & c_{0,1}^l & \dots & c_{0,n-1}^l \\ c_{1,0}^l & c_{1,1}^l & \dots & c_{1,n-1}^l \\ \vdots & \vdots & \vdots & \vdots \\ c_{n-1,0}^l & c_{n-1,1}^l & \dots & c_{n-1,n-1}^l \end{bmatrix}$$

• The shortest path problem has been formulated as a sequence of *r* matrix-vector products.

Parallel Shortest-Path

- We can parallelize this algorithm using the parallel algorithms for the matrix-vector product.
- $\Theta(n)$ processing elements can compute each vector C' in time $\Theta(n)$ and solve the entire problem in time $\Theta(rn)$.
- In many instances of this problem, the matrix *M* may be sparse. For such problems, it is highly desirable to use sparse matrix techniques.

- We are given a knapsack of capacity c and a set of n objects numbered 1,2,...,n. Each object i has weight w_i and profit p_i.
- Let $v = [v_1, v_2, ..., v_n]$ be a solution vector in which $v_i = 0$ if object *i* is not in the knapsack, and $v_i = 1$ if it is in the knapsack.
- The goal is to find a subset of objects to put into the knapsack so that

$$\sum_{i=1}^{n} w_i v_i \le c$$

(that is, the objects fit into the knapsack) and

$$\sum_{i=1}^{n} p_i v_i$$

is maximized (that is, the profit is maximized).

- The naive method is to consider all 2ⁿ possible subsets of the n objects and choose the one that fits into the knapsack and maximizes the profit.
- Let *F[i,x]* be the maximum profit for a knapsack of capacity *x* using only objects {1,2,...,i}. The DP formulation is:

$$F[i,x] = \left\{egin{array}{ccc} 0 & x \geq 0, i=0\ -\infty & x < 0, i=0\ \max\{F[i-1,x], (F[i-1,x-w_i]+p_i)\} & 1 \leq i \leq n \end{array}
ight.$$

- Construct a table *F* of size *n x c* in row-major order.
- Filling an entry in a row requires two entries from the previous row: one from the same column and one from the column offset by the weight of the object corresponding to the row.
- Computing each entry takes constant time; the sequential run time of this algorithm is $\Theta(nc)$.
- The formulation is serial-monadic.

Table F

n Fij i 2 1 Weights -1 j-wi j c-1 С Processors Pi-1 **P0** Pi-wi-1 Pc-2 Pc-1

Computing entries of table *F* for the 0/1 knapsack problem. The computation of entry *F*[*i*,*j*] requires communication with processing elements containing entries *F*[*i*-1,*j*] and *F*[*i*-1,*j*-*w*_{*i*}].

- Using c processors in a PRAM, we can derive a simple parallel algorithm that runs in O(n) time by partitioning the columns across processors.
- In a distributed memory machine, in the *jth* iteration, for computing *F[j,r]* at processing element *P_{r-1}*, *F[j-1,r]* is available locally but *F[j-1,r-w_j]* must be fetched.
- The communication operation is a circular shift and the time is given by $(t_s + t_w) \log c$. The total time is therefore $t_c + (t_s + t_w) \log c$.
- Across all *n* iterations (rows), the parallel time is O(n log c). Note that this is not cost optimal.

- Using *p*-processing elements, each processing element computes *c/p* elements of the table in each iteration.
- The corresponding shift operation takes time $(2t_s + t_w c/p)$, since the data block may be partitioned across two processors, but the total volume of data is c/p.
- The corresponding parallel time is $n(t_c c/p + 2t_s + t_w c/p)$, or O(nc/p) (which is cost-optimal).
- Note that there is an upper bound on the efficiency of this formulation.

Nonserial Monadic DP Formulations: Longest-Common-Subsequence

- Given a sequence $A = \langle a_1, a_2, ..., a_n \rangle$, a subsequence of A can be formed by deleting some entries from A.
- Given two sequences $A = \langle a_1, a_2, ..., a_n \rangle$ and $B = \langle b_1, b_2, ..., b_m \rangle$, find the longest sequence that is a subsequence of both *A* and *B*.
- If A = <*c*,*a*,*d*,*b*,*r*,*z*> and B = <*a*,*s*,*b*,*z*>, the longest common subsequence of A and B is <*a*,*b*,*z*>.

Longest-Common-Subsequence Problem

- Let *F[i,j]* denote the length of the longest common subsequence of the first *i* elements of *A* and the first *j* elements of *B*. The objective of the LCS problem is to find *F[n,m]*.
- We can write:

$$F[i,j] = \left\{ egin{array}{ll} 0 & ext{if } i=0 ext{ or } j=0 \ F[i-1,j-1]+1 & ext{if } i,j>0 ext{ and } x_i=y_j \ \max{\{F[i,j-1],F[i-1,j]\}} & ext{if } i,j>0 ext{ and } x_i
eq y_j \end{array}
ight.$$

Longest-Common-Subsequence Problem

- The algorithm computes the two-dimensional F table in a row- or column-major fashion. The complexity is $\Theta(nm)$.
- Treating nodes along a diagonal as belonging to one level, each node depends on two subproblems at the preceding level and one subproblem two levels prior.
- This DP formulation is nonserial monadic.

Longest-Common-Subsequence Problem



(a) Computing entries of table for the longest-commonsubsequence problem. Computation proceeds along the dotted diagonal lines. (b) Mapping elements of the table to processing elements.

Longest-Common-Subsequence: Example

• Consider the LCS of two amino-acid sequences H E A G A W G H E E and P A W H E A E. For the interested reader, the names of the corresponding amino-acids are A: Alanine, E: Glutamic acid, G: Glycine, H: Histidine, P: Proline, and W: Tryptophan.

		Η	Ε	Α	G	Α	W	G	Η	Ε	Ε
	0	0	0	0	0	0	0	0	0	0	0
Р	0	0	0	0	0	0	0	0	0	0	0
Α	0	0	0	1	1	1	1	1	1	1	1
W	0	0	0	1	1	1	2	2	2	2	2
н	0	1	1	1	1	1	2	2	3	3	3
E	0	1	2	2	2	2	2	2	3	4	4
A	0	1	2	3	3	3	3	3	3	4	4
Е	0	1	2	3	3	3	3	3	3	4	5

• The *F* table for computing the LCS of the sequences. The LCS is A W H E E.

Parallel Longest-Common-Subsequence

- Table entries are computed in a diagonal sweep from the top-left to the bottom-right corner.
- Using *n* processors in a PRAM, each entry in a diagonal can be computed in constant time.
- For two sequences of length *n*, there are 2*n*-1 diagonals.
- The parallel run time is Θ(n) and the algorithm is costoptimal.

Parallel Longest-Common-Subsequence

- Consider a (logical) linear array of processors. Processing element P_i is responsible for the $(i+1)^{th}$ column of the table.
- To compute F[i,j], processing element P_{j-1} may need either F[i-1,j-1] or F[i,j-1] from the processing element to its left. This communication takes time $t_s + t_w$.
- The computation takes constant time (t_c) .
- We have:

$$T_P = (2n-1)(t_s + t_w + t_c).$$

- Note that this formulation is cost-optimal, however, its efficiency is upper-bounded by 0.5!
- Can you think of how to fix this?

Serial Polyadic DP Formulation: Floyd's All-Pairs Shortest Path

- Given weighted graph G(V,E), Floyd's algorithm determines the cost d_{ij} of the shortest path between each pair of nodes in V.
- Let d^k_i be the minimum cost of a path from node i to node j, using only nodes v₀, v₁, ..., v_{k-1}.
- We have:

$$d_{i,j}^k = \left\{ egin{array}{cl} c_{i,j} & k = 0 \ \min \left\{ d_{i,j}^{k-1}, (d_{i,k}^{k-1} + d_{k,j}^{k-1})
ight\} & 0 \leq k \leq n-1 \end{array}
ight.$$

• Each iteration requires time $\Theta(n^2)$ and the overall run time of the sequential algorithm is $\Theta(n^3)$.

Serial Polyadic DP Formulation: Floyd's All-Pairs Shortest Path

- A PRAM formulation of this algorithm uses n² processors in a logical 2D mesh. Processor P_{i,j} computes the value of d^k_{i,j} for k=1,2,...,n in constant time.
- The parallel runtime is $\Theta(n)$ and it is cost-optimal.
- The algorithm can easily be adapted to practical architectures, as discussed in our treatment of Graph Algorithms.

Nonserial Polyadic DP Formulation: Optimal Matrix-Parenthesization Problem

- When multiplying a sequence of matrices, the order of multiplication significantly impacts operation count.
- Let C[i,j] be the optimal cost of multiplying the matrices A_i,...A_j.
- The chain of matrices can be expressed as a product of two smaller chains, A_i, A_{i+1}, ..., A_k and A_{k+1}, ..., A_j.
- The chain $A_{i}, A_{i+1}, \dots, A_k$ results in a matrix of dimensions $r_{i-1} \times r_k$, and the chain A_{k+1}, \dots, A_j results in a matrix of dimensions $r_k \times r_j$.
- The cost of multiplying these two matrices is $r_{i-1}r_kr_i$.

• We have:

$$C[i,j] = \begin{cases} \min_{\substack{i \le k < j \\ 0}} \{C[i,k] + C[k+1,j] + r_{i-1}r_kr_j\} & 1 \le i < j \le n \\ 0 & j = i, 0 < i \le n \end{cases}$$



A nonserial polyadic DP formulation for finding an optimal matrix parenthesization for a chain of four matrices. A square node represents the optimal cost of multiplying a matrix chain. A circle node represents a possible parenthesization.

- The goal of finding *C*[1,*n*] is accomplished in a bottom-up fashion.
- Visualize this by thinking of filling in the C table diagonally. Entries in diagonal / corresponds to the cost of multiplying matrix chains of length I+1.
- The value of C[i,j] is computed as $min\{C[i,k] + C[k+1,j] + r_{i-1}r_kr_j\}$, where k can take values from i to j-1.
- Computing C[i,j] requires that we evaluate (j-i) terms and select their minimum.
- The computation of each term takes time t_c, and the computation of C[i,j] takes time (j-i)t_c. Each entry in diagonal I can be computed in time It_c.

- The algorithm computes (n-1) chains of length two. This takes time (n-1)t_c; computing n-2 chains of length three takes time (n-2)t_c. In the final step, the algorithm computes one chain of length n in time (n-1)t_c.
- It follows that the serial time is $\Theta(n^3)$.



The diagonal order of computation for the optimal matrixparenthesization problem.

Parallel Optimal Matrix-Parenthesization Problem

- Consider a logical ring of processors. In step *I*, each processor computes a single element belonging to the *I*th diagonal.
- On computing the assigned value of the element in table *C*, each processor sends its value to all other processors using an all-to-all broadcast.
- The next value can then be computed locally.
- The total time required to compute the entries along diagonal *I* is $lt_c+t_s log n+t_w(n-1)$.
- The corresponding parallel time is given by:

$$T_P = \sum_{l=1}^{n-1} (lt_c + t_s \log n + t_w(n-1)),$$

$$=\frac{(n-1)(n)}{2}t_c+t_s(n-1)\log n+t_w(n-1)^2.$$

Parallel Optimal Matrix-Parenthesization Problem

- When using *p* (*<n*) processors, each processor stores *n/p* nodes.
- The time taken for all-to-all broadcast of *n/p* words is

$$t_s\log p + t_w n(p-1)/p pprox t_s\log p + t_w n$$

and the time to compute n/p entries of the table in the l^{th} diagonal is $lt_c n/p$. $T_P = \sum_{l=1}^{n-1} (lt_c n/p + t_s \log p + t_w n),$

$$=\frac{n^2(n-1)}{2p}t_c+t_s(n-1)\log p+t_wn(n-1).$$

 $T_P = \Theta(n^3/p) + \Theta(n^2)$.

 This formulation can be improved to use up to n(n+1)/2 processors using pipelining.

Discussion of Parallel Dynamic Programming Algorithms

- By representing computation as a graph, we identify three sources of parallelism: parallelism within nodes, parallelism across nodes at a level, and pipelining nodes across multiple levels. The first two are available in serial formulations and the third one in non-serial formulations.
- Data locality is critical for performance. Different DP formulations, by the very nature of the problem instance, have different degrees of locality.