

Backtracking Branch & Bound

Backtracking - It is a general algo technique

Principle for solving computational problems incrementally by trying out various possibility and abandoning those that fail to meet the problem's constraint. The idea is to explore the solution space systematically backtracking when necessary, and try alternative choices.

> Principles of Backtracking -

- 1) The solution is built incrementally by making a series of choices at each step decision is made.
- 2) If a chosen path leads to a state that can't possibly be part of valid soln' the algo backtracks and abandons the path (Prune).
- 3) The soln' Space is explored Systematically, often using recursion. Algo explores one branch of decision tree at a time.

Q)

> Control Abstraction -

- 1) In backtracking, the soln' is defined as an tuple $x = (x_1, x_2, \dots, x_n)$ where $x_i = 0$ or x_i is chosen from set of finite components S_i . If component $x_i = 1$ else set it to 0. The backtracking approach tries to find vector x such that it maximizes or minimizes certain criterium $F(x)$.

B is Bounding Function.

Recursive approach - Algorithm.

1. Algorithm Backtrack(K)

```

{ For each  $x[K] \in T(x[1], x[2], \dots, x[K-1])$  do
  { if ( $B_K(x[1], \dots, x[K]) \neq 0$ ) then
    { if ( $(x[1], x[2], \dots, x[K])$  is a path to answer node)
      then write ( $x[1:K]$ );
      if ( $K < n$ ) then Backtrack ( $K+1$ );
    }
  }
}

```

3

3

9

> Complexity Analysis -

- > Time to compute next candidate solution $x[i]$
- > Number of candidate soln $x[i]$ fulfilling the explicit constraints.
- > Time taken by bounding func n B_i to check promising candidate soln
- > No. of candidate soln $x[i]$ fulfilling satisfying the bounding func n B_i .
- > For a problem instance of size n , if the backtracking algo produces a soln space of 2^n or $n!$ nodes, then it has worst case running times $O(p(n)2^n)$ or $O(q(n)n!)$ respectively where $p(n)$ & $q(n)$ are polynomials in n .

* Basic Backtracking terminologies -

> Solution vector - The desired soln x to a problem instance p of input size n is expressed as vector of candidate soln x_i that are selected from finite set of possible soln S . Thus a soln can be represented as an \mathbb{Z}^n -tuple (x_1, x_2, \dots, x_n) & partial soln (x_1, x_2, \dots, x_i)

> Constraints - This are the rules to confine the soln vector (x_1, x_2, \dots, x_n) . They determine the values of candidate soln and their relationship with each other.

2 types

> Implicit & > Explicit.

> Implicit Constraints -

→ These are the rules that identify the tuples in the solution space S that satisfy the specified criterion f_{cone} of problem instance P .

2) example N Queen.

? Explicit Constraints - This are the rules by which all candidate solutions are restricted to take on values from a specified set in a problem instance P .

~~A~~ Iterative backtracking Approach.

Algorithm Backtrack()

$k := 1$

 while $k \neq 0$ do

 if ($\text{untried } x[k] \in T[x[1], x[2], \dots, x[k-1]]$) AND

 ($B_k(x[1], x[2], \dots, x[k])$ is path to answer node) then

 Print $x[1], x[2], \dots, x[k]$ // soln?

$k \leftarrow k + 1$ // consider next candidate

 else

$k \leftarrow k - 1$ // Backtrack to recent Node

 end

 end.

N-Queen - The N Queen is a classical combinatorial problem that asks for all arrangements of N queen on $N \times N$ chess board, such that no two queens threaten each other.

(consider an $n \times n$ chess board on which)

- 1) Initially chess board is empty
- 2) Reasons why backtracking is needed explain in own words first without Backtrack then with backtrack).

Pseudocode -

Algorithm Queen (n)

For Column $\leftarrow 1$ to n do

{ if (place (row, column)) then
 { board [row] [column] // no conflict, place queen
 if (row = n) then // dead end
 print board (n) // printing board
 else // try next queen with next pos?
 Queen (row + 1, n)
 } }

Row by Row each queen
 is placed by satisfying
 constraints.

Algorithm place (row, column)

for i $\leftarrow 1$ to row - 1 do

{ if (board [i] = column) then
 return 0. // check in same column.
 else if (abs (board [i] - column) = abs (i - row)) then
 return 0 // check diagonally

/ no conflicts then queen can be placed.

return 1.

Graph Colouring Problem - Consider a Graph $G(V, E)$ V is set of vertices & E is set of edges. Then the graph coloring problem asks to assign m colors to the vertices of G such that no two non-neighbouring vertices have a similar color.

Pseudocode -

Algorithm ColorGraph(G, k):

 colors = $\{0\}^* | V|$

 colorVertex($G, 0, k, \text{colors}$)

Algorithm colorVertex($G, \text{vertex}, k, \text{colors}$):

 if vertex is to $|V|$ then

 return True. // All vertices are successfully colored.

 for color from 1 to k do

 if isSafe($G, \text{vertex}, \text{color}, \text{colors}$) is true then

 // Assign color to vertex

 colors[vertex] = color

 if colorVertex($G, \text{vertex}+1, k, \text{colors}$) is true then

 return True. // Coloring successful.

 // If coloring fails

 colors[vertex] = 0

 // If no color is suitable return False.

 return False.

Algorithm isSafe($G, \text{vertex}, \text{color}, \text{colors}$):

 for adjacent vertex in neighbors(G, vertex) do

 if colors[adjacent vertex] is equal to color then

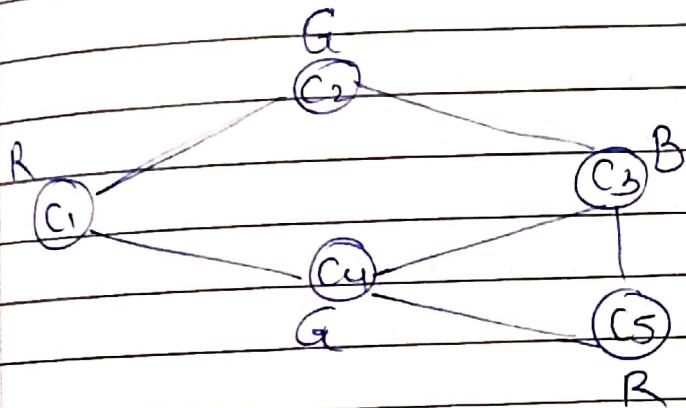
 return False // color conflict

 // If no conflict are found, color is safe.

 return True

	C ₁	C ₂	C ₃	C ₄	C ₅
C ₁	0	1	0	1	0
C ₂	1	0	1	0	0
C ₃	0	1	0	1	1
C ₄	1	0	1	0	1
C ₅	0	0	1	1	0

Color = R G B.



~~* Branch & Bound~~ - It systematically explores the Soluⁿ Space of a problem dividing it into smaller subproblems, solving them & bounding the Soluⁿ Space to eliminate Subproblems that can't lead to an optimal Soluⁿ

> Characteristics -

1) Divide & Conquer -

2) Exploration of Soluⁿ Space

3) Bounding Techniques - It involves establishing upper bound & lower bounds. Subproblems that cannot lead to optimal Soluⁿ then they are pruned.

4) Priority Queue = Algo uses priority Queue

5) Branching & Pruning

6) Terminal criteria.

> General Algo -

Algorithm Branch & Bound (S)

E \leftarrow createNode

while true do

 if E is the leaf node then

 print "Display path from E"

 return

 end

 Expand (E)

 if size (Heap) = = 0 then

 print "Soln not found"

 return

 end

 E \leftarrow deque (H)

end

Algorithm Expand (E)

Generate all

possible child

nodes of E

Compute the cost

of each child C.

Insert each child

C in Heap.

Search Techniques -

1) LC Search (Least Cost Search) - It uses heuristic cost funcⁿ to compute the bound values at each node. Nodes are added to the list of live nodes as soon as they get generated. The node with least value of a cost function is selected as a next E-node.

2) BFS FIFO - It maintains the list of live nodes in first in first out order in a queue.

3) DFS LIFO - Uses Stack

4) Control Abstraction for FIFO / BFS

Algorithm BB-FIFO()

{ if T is the answer node Then

{ mm (cost[T], ub[T] + E)

 write T

 return ;

}

E := T; // E-node

Initialize empty Queue of live nodes.

repeat

{ For (each child K of E)

{ if (K is an answer state)

{ write path from K to T;

 return ;

}

 Append Q(K);

(K → Parent) := E;

}

if (there are no more live nodes)

{ write ("No answer state");

 return ;

E := Delete_Q() // Delete live node at the front & return it

{ until (FALSE);

}

> Control Abstract for LIFO / DFS

~~TypeDef Struct.~~

~~{ ListNode - T* Next;~~

~~ListNode - T* parent;~~

~~Float cost~~

~~} ListNode - T is the state space tree~~

Algorithm BB-LIFO()

if T is the answer node Then

$u \leftarrow \min \{ \text{cost}(T), \text{ub}(T) + E \}$

Prmt(T)

return

end end

$E \rightarrow E \pm$

$E \leftarrow T$

while true do

For each child x of T do

if x is the answer node then

Prmt "display path from x to root"

return

end

PUSH(x) //push new node on top of stack

$x \rightarrow \text{parent} := E$

if x is answer node and $\text{cost}(x) \leq \text{ub}$ then

$u \leftarrow \min \{ \text{cost}(T), \text{ub}(T) + E \}$

end

if no more live nodes then

Prmt "No live nodes"

Prmt "Minimucast : ub"

return

end

$E \leftarrow \text{POPC}$

end.

Control abstract for LC Search - LC Search uses heuristic function to assign the ranks to live nodes.
It estimates the extra computational work (cost) to reach an answer state from current live node.

The minimum number of nodes required to be generated in subtree of k to reach an answer state.

$$f(k) = F(h(k)) + g^*(k) \text{ where } F(\cdot) \text{ is an increasing func.}$$

Algorithm BB-LC_Search()

if T is the answer node then

 print(T)

 return

end

$E \leftarrow T$

while (True) do

 For each child X of E do

 if X is the answer node then

 print "Display path from X to root"

 return

 end

 INSERT(X) // Insert in to list

$X \rightarrow \text{parent} := E$

end

if no more live nodes then

 print "soln Found"

 return

end.

$iE \leftarrow \text{least_cost}()$ // Find next least node from list

end

Amortized Analysis

- > Amortized analysis is a technique used in CS to analyze the average case time complexity of algorithms that perform a sequence of operations, where some operan may be expensive than others. The idea is to spread the cost of these expensive operations over multiple operan. So that the average cost of each insertion becomes constant or less.
- > Amortized analysis is useful in designing algorithms for data structures such as dynamic arrays, Priority queues. Let us consider ex. of simple hash table. How do we decide on Table size? The idea is to increase the size of table whenever it is full and copy content in new table & then forget older one.
- > Worst case time is $O(n)$ of inserⁿ is $O(n)$ worst case of n inserts is $n \cdot O(n)$ which is $O(n^2)$.

Techniques of Amortized Analysis -

- 1) Aggregate Method 2) Accounting 3) Potential method

> **Aggregate Method** - This technique involves analyzing a sequence of operations as a whole rather than focusing on individual operan. The total cost of a sequence of operan is divided by the number of operations to obtain the avg cost.

Example 1 : Stack operation.

Consider a stack data structure it performs 2 operan

1) PUSH(x, p) 2) POP(x)

∴ Total cost of a sequence that contains n PUSH & POP operan = $\Theta(n)$

Let us consider MultiPOP(x, m) which pops the topmost m objects of stack x.

Let us consider an empty stack x and n PUSH, POP & MULTIPOP operan are executed on stack

As there are n PUSH operatⁿ executed on the Stack & stack size will be n so the worst case cost of a single MULTIPOP operatⁿ will be $O(n)$ operatⁿ worst case cost.

PUSH (x, p)	$O(1)$
POP (x)	$O(1)$
MULTIPOP (x, m)	$O(n)$

Worst case of any stack operatⁿ is $O(n)$ then n operatⁿ on stack will be $O(n^2)$. We can pop at most n object PUSH onto the stack. So the total cost of any sequence of n PUSH, POP & MULTIPOP operatⁿ on stack will be $O(n)$ for $\forall n$. Thus we get Avg cost of each stack operatⁿ in sequence of n PUSH, POP, MULTIPOP = $\frac{O(n)}{n} = O(1)$

This is Amortized Cost.

> Accounting Method - When the amortized cost of operatⁿ is assigned to the particular object and some are assigned lesser amortized cost than actual cost.

> When Amortized cost is larger than its actual cost, the excess amount is assigned to the particular object in data structure as its credit. This credit is consumed for other operation whose amortized cost is lesser than its actual cost.

> Stack Example.

Stack operations with PUSH(x,p) POP(x) MULTIPOP(x,m)		
Opn	Actual cost	Amortized cost
PUSH(x,p)	1	2
POP(x)	1	0
MULTIPOP(x,m)	min(Stack size, m)	0

> for n PUSH operation the amortized cost will be 2n whereas the actual cost of n PUSH opern will be n so total credit = $2n - n = n$ will be assigned to stack object.

> Actual cost of single POP(x) is 1 and credit of stack object after n PUSH opern is n. Therefore. POP(x) Separately or MULTIPOP(x) can consume the credit (n). Thus for any sequence of n PUSH POP & MULTIPOP opern AS in This example total amortized cost = $O(n)$ the total actual cost of n stack opern is $O(n)$.

> Potential Function Method - This is similar to Accounting method but rather than credit, it uses potential or energy to pay the cost of operation.

Consider that n operations are carried out on a data structure with initial state S_0 , let S_i be the state of whole data structure after performing on i^{th} opern

> The function ϕ is selected such that it maps each state of data structure S_i to a real no. $\phi(S_i)$ ϕ is known as "Potential function" and $\phi(S_i)$ is called as "Potential Energy"

The amortized cost of i^{th} operan is \hat{a}_i

$$\hat{a}_i = a_i + (\phi(s_i) - \phi(s_{i-1}))$$

Total Amortized cost = $\sum_{i=1}^n \hat{a}_i$

$$\sum_{i=1}^n \hat{a}_i = \sum_{i=1}^n (a_i + \phi(s_i) - \phi(s_{i-1}))$$

> Stack example - PUSH POP MULTIPOP operations

Considered Potential Function ϕ is defined on the stack size. Initially for empty stack $\phi(s_0) = 0$

s_i resulting after the i^{th} operan will have potential $\phi(s_i) \geq 0$

1) PUSH(x, k) - As it pushes on a single element onto stack of size k the stack size will increase to $k+1$. Actual cost of 1 PUSH operan is 1

$$\begin{aligned}\hat{a}_i &= a_i + [\phi(s_i) - \phi(s_{i-1})] \\ &= 1 + [(k+1) - k] = 1\end{aligned}$$

2) POP(x) - As it pops stack size decreases to $k-1$

$$\begin{aligned}\hat{a}_i &= a_i + [\phi(s_i) - \phi(s_{i-1})] \\ &= 1 + [(k-1) - k] = 0\end{aligned}$$

3) MULTIPOP(x, m) - $2 = \min(k, m)$ The actual cost of operan is 2

$$\begin{aligned}\hat{a}_i &= a_i + [\phi(s_i) - \phi(s_{i-1})] \\ &= 2 + [(k-2) - k] = 0.\end{aligned}$$

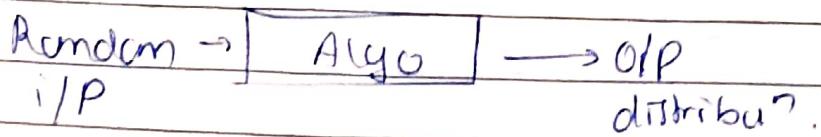
Since Amortized cost of each PUSH, POP, MULTIPOP operan is $O(n)$ $O(1)$ the total Amortized cost of sequence of n PUSH POP MULTIPOP operan is $O(n)$

★ Tractable Problems - Feasibility of an algorithm to complete its exec'n in reasonable time. If they get solved in polynomial time using deterministic algorithms Time complexities $O(n^2)$ ($\Theta(n^2)$), $O(1)$ example - Given an array of n elements, the problem is to arrange elements in ascending or descending order Algos like QuickSort, mergeSort & HeapSort
 Linear Search - $O(n)$ binary search $O(\log n)$

★ Non tractable problems - The infeasibility of an algorithm to complete its exec'n in reasonable time Pro if problems cannot be solved in polynomial time using deterministic algorithm Time complexities like. $O(2^n)$, $O(n^n)$, $O(n!)$
 ex 01 Knapsack $O(2^{n+2})$, TSP $O(n^{2^n})$

★ Randomized Algorithms - An Algorithm that makes some random choices during its exec'n depending on random numbers it's for its operations
 Randomized algo use randomness in their decision making process. This randomness can come from various sources.

? It is also known as "probabilistic algorithm" as input is supposed to be from a probability distribution



? The randomized algo avoid worst case behaviour of classic deterministic algo

- > These algo are non-deterministic. These algo can make random but most of time correct decision.
- > These algo may operate at different efficiencies in different runs for same instance of problem.
- > classes of randomized algo
 - > Las Vegas algo - This algo always computes the same result for same input. The running time of these algo vary with the o/p randomized. The ~~new~~ running time may vary.
 - > Monte Carlo - Always computes diff. results in different runs for same input. Produces output with some probability for fixed input. They do not show much variation.

*** Approximation Algo** - it is an algo which efficiently finds a solution to an optimization problem that is guaranteed to be close to the optimal solution. Finding exact optimal soln may be computationally infeasible.

> Finding an optimal solution for NP-complete is not polynomial bounded. Approximation algo apply some heuristic to cut off the further computation in each step.

> The quality of solⁿ produced by an approximation is measured by its approximation ratio.

$$\text{Approx Ratio } P = \frac{\text{cost of Approximate Sol}^n}{\text{optimal cost}}$$

> The approx ratio provides a performance guarantee, ensuring that the solⁿ produced by the algo is within a certain factor of optimal solⁿ.

Ex -> Greed Algo \Rightarrow Heuristic Methods

Embedded Algo - A computer hardware with software embedded in it is called an Embedded Sys. It's a reliable, microcontroller or microprocessor based software driven realtime control System.

Characteristics -

- 1) Single Functioned -
 - 2) Tightly constrained - Size, cost, memory, power
 - 3) Realtime
 - 4) Microprocessor based.
 - 5) Memory -
- > The algo implemented on microcontroller or microprocessor based system is embedded Algo. They solve specialized oprcsⁿ in real time embedded algo face challenges like. limited memory, low power consumeⁿ, deadline based Scheduling, less cost.

Multithreaded & Distributed Algo.

A thread is a dispatchable portion of a task within a process. Each thread executes its code sequentially. The algorithm that permit the concurrent exec'n of multiple threads of a program on multiple core in system with shared memory are known as multithreaded algo.

→ Static multithreaded algo - The no. of threads and their exec'n behaviour is determined before the program starts running. Programmer explicitly specifies the no. of threads and the structure of parallelism is fixed at compile time.

→ Dynamic multithreaded algo - The no. of threads and their exec'n behaviour can change during exec'n. Threads are created managed and terminated dynamically based on program's runtime behaviour & workload.

→ Different keywords -

1) parallel - Used with the loop construct like for loops to mention that all iterat' com concurrently executed.

2) spawn - Used to provide the nested parallelism.

In a general func'n call, the parent doesn't resume until its child returns, the parent & child executed concurrently.

3) Sync - A procedure com safely return values of its spawned children. The procedure is suspended until all its children have executed completely. Then the procedure resumes the instruc'n following Sync Step.

* Performance measure -

1) Work - Total time required to execute the entire computer on one process, so the work represents the sum of the time taken by each thread. An ideal parallel computer with P processors can perform up to P units of work in one time step. It can do $P \times T_p$ work in T_p time. Since work is - Total work is $T_1 \cdot P \times T_p \geq T$, Hence work done is $T_p \geq \frac{T}{P}$.

2) Span - It is the longest time it takes to execute threads along any path of the Computational DAG (Directed Acyclic Graph).

An ideal parallel computer with P processors can't run faster than a machine with an infinite no. of processor. On other hand, can imitate a P processor machine by employing only $\frac{1}{P}$ of its processor. $P \cdot T_p \geq T$ which is Span done.

3) Speedup & Parallelism - Speedup of computer on P processors is $\frac{T_1}{T_P}$. Using work done, we can

conclude that an ideal parallel computer can not have any more speedup than no. of processors. The parallelism of multithreaded computer is $\frac{T_1}{T_{avg}}$. So this ratio means the average amount of T_{avg} work that can be performed for each step of parallel execution.

4) Slackness - it is parallelism by P ratio

$$\frac{\left(\frac{T_1}{T_{avg}}\right)}{P}$$

Analyzing multithreaded Algos.

> Analyzing the work - The work done by a multithreaded algorithm is analyzed by ignoring the parallel constructs and concentrating on the analysis of sequential algo

$$T_1(n) = T(n) = T(n-1) + T(n-2) + \Theta(1)$$

$$T(n) = \Theta(F_n)$$

where F_n grows exponentially in $F_n = \phi^n$ where $\phi = \alpha$

$$\text{Golden ratio} = \left(\frac{1+\sqrt{5}}{2}\right)$$

Analyzing Spans -

For serial composition of subcomputations - If a set of subcomputations are in series then the span of their composition is obtained by the sum of the spans of those sub-computations

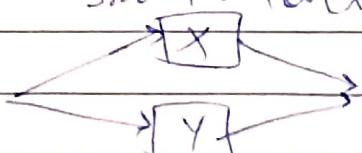
For parallel composition - If a set of parallel subcomputations are in parallel, then the span of their composition is the maximum of the spans of those sub computations

The work & span analysis.



work $T_1(X \cup Y) = T_1(X) + T_1(Y)$ Series,

span $= T_{\infty}(X \cup Y) = T_{\infty}(X) + T_{\infty}(Y)$



work $= T_1(X \cup Y) = T_1(X) + T_1(Y)$

span $= T_{\infty}(X \cup Y) = \max(T_{\infty}(X), T_{\infty}(Y))$.

> Analyzing Parallel Loops - At multithread Algos

use keyword `parallel` with the loop constructs to define parallel execution of iteration. To design parallel algo to multiply $n \times n$ matrix $X = x_i$ with vector y_i of size $n \times 1$

$$z_i = \sum_{j=1}^n x_{ij} y_j$$

Algorithm `Parallel_Multi(x, y)`:

`parallel for i ← 1 to n do`

$$z_i, z_j \leftarrow 0$$

`end.`

`parallel for i ← 1 to n do`

`for j ← 1 to n do`

$$z_i \leftarrow z_i + (x_{ij} * y_j)$$

`end.`

`end.`

Keyword `parallel for` indicates that iteration of the loop may run concurrently.

Algorithm `matvec_MUL-Loop(x, y, z, n, i, i')`

`if i < i' then`

`for j ← 1 to n do`

$$z_i \leftarrow z_i + (x_{ij} * y_j)$$

`end`

`else mid ← (i + j) / 2`

`Spawn MATVEC-MUL-Loop(x, y, z, n, i, mid)`

`MATVEC-MUL-Loop(x, y, z, n, mid+1, i')`

`Sync`

`end`

The algo recursively Spans First half of the loop which runs parallel with Second half

work done by parallel matrix is simple case of matrix

$$T(n) = O(n^2)$$

$$\text{Span is } T_{\infty}(n) = \Theta(\log n).$$

> Race condition - Algos are said to be deterministic if it always does the same work for the same input, irrespective of how the instrucⁿ are scheduled on a multicore machine. When 2 or more threads access shared data concurrently, & at least one of them modifies the data.

Race Condition

$$x \leftarrow 10$$

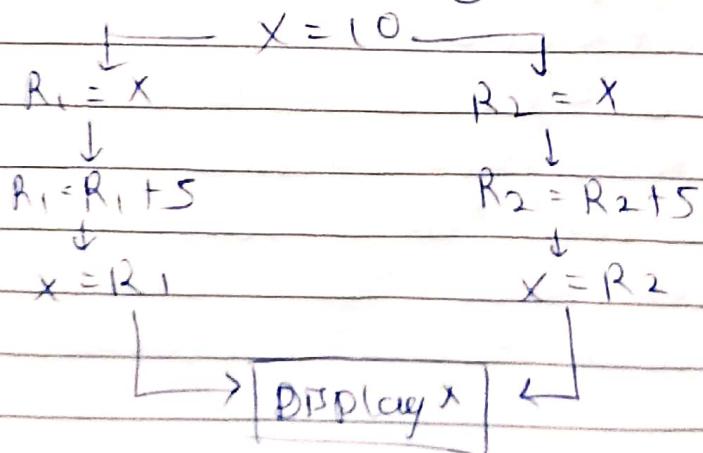
parallel for $i \leftarrow 1$ to 2 do

$$x \leftarrow x + 5$$

end.

After initializing x to 10 the parallel loop creates two strands each performing addⁿ to x .

The expected outcome of code is 20 but algo might return 15. Two strands run in parallel on 2 processors. Both processors loads x , update x , and write back to the memory.



Multithreaded Matrix Multiplication -

P-matrix-Multiply(A, B)

$n = A \times B \rightarrow n^2$

If C is a new $n \times n$ matrix.

parallel For $i=1$ to n do

parallel For $j=1$ to n do

$C_{ij} = 0$

For $k=1$ to n do

$C_{ij} = C_{ij} + a_{ik} \times b_{kj}$

end

end

return C

end.

In this algo the longest path is when spawning the outer and inner parallel loop execn and then the n execn for innermost for loop.

So the Spm of this algo is $T_{\text{sp}}(n) = O(n)$

$$\frac{T_1(n)}{T_{\text{sp}}(n)} = \frac{O(n^2)}{O(n^2)} = 1$$

Multithreaded Merge Sort -

~~Merge-Sort(A, p, r)~~ At

~~if P < r then~~

~~q = (P+r)~~

Algorithm merge-Sort(x, first, last)

{ if (first < last)

{ mid := (first + last) / 2

Spawn mergesort(x, first, mid);

mergesort(x, mid+1, last);

Syn C; }

merge-results(x, first, mid, last);

}

Since our algo Merge-Result() remains a sequential algo, its work and Spcm are $\Theta(n)$ where n is input size.

Work of merge-sort() is $T(n) =$

$$T(n) = 2 \left(2 T(n/2) + \Theta(n) \right) = \Theta(n \log n)$$

and the Spcm $T_{\text{sp}}(n)$ of merge-sort()

$$T_{\text{sp}}(n) = T_{\text{sp}}(n/2) + \Theta(n) = \Theta(n).$$

$$\text{Parallelism} = \frac{T_{\text{sp}}(n)}{T_{\text{sp}}(n)} = \frac{\Theta(n \log n)}{\Theta(n)} = \Theta(\log n).$$

~~Distributed BFS -~~

A parallel Algo that facilitates concurrent execution in distributed computing environment.

A BFS algo is applied to traverse or search a graph or tree data structure. distributed BFS performs over BFS over distributed network

- > In 1-D partitioning a graph is partitioned such that each node & edges originating from it are possessed by one processor
- > P is the no. of processors. Edges emanating from vertex v is a list of vertex index in row v of adjacency matrix. process -

> 1) Create a set of Frontier vertices (F) owned by processor

2) Edge lists of F are merged to form a set of neighbor vertices N

3) Send msg to processes that own vertices in set N and add these vertices in their next F

4) Receive set N from other processes

5) Merge to form final set N

6) Update level.

A Rabin Karp algo - It is a String searching algo

- It computes hash value of the pattern and hash value of overlapping such strings of text.
- Compare hash values. If the hash value matches the text is detected in pattern
- Rolling Hash - To efficiently update the hash value when moving from one substring to the next.

Pseudocode -

Function rabin-karp (text, pattern) :

n = length of text

m = length of pattern

P-hash = hash (pattern)

text-hash = hash (text[0:m])

for i from 0 to n-m :

if text-hash == P-hash :

if text[i]

if text[i : itm] == pattern

print ("Found")

if i < n-m :

text-hash = rehash (text-hash, text[i],

text[itm])

Function hash (s) :

result = 0

For char in s

result = (result * base + ord(char)) % prime

return result.