A* Path-Finding with MPI

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Problem Description

 We want to find the shortest path between two points on a connected graph

 We have a known start point, end point, and know the cost to move into any adjacent node from every point



A* Algorithm

• Similar to other search algorithms like Dykstra's or breadth-first

• Uses a heuristic function to make decisions on traversal of graph

- The total cost from start to finish is modeled by F(n) = G(n) + H(n)
 - F(n): the total cost
 - G(n): cost to move into node we are on
 - H(n): Heuristic function estimating remaining cost to goal

A* Algorithm

- When(if) a path is found from start to finish we want this F cost to be minimal
- This is where the heuristic function is used
- By estimating the remaining distance to the goal we can make an educated guess on where to check next



Optimal path found

A* Algorithm - Heuristic Function

- There are a wide variety of metrics to consider for the heuristic
 - Optimality of solution
 - Completeness of solution set
 - Accuracy
 - Running time
- Must consider what we need and tradeoffs associated
- Common choices
 - Manhattan: x,y vector distance to goal
 - Euclidean : straight line distance to goal
 - Chebyshev: max of either x or y distance to goal

A* Algorithm - Heuristic Function

• Heuristic function must not overestimate remaining distance to be considered admissible





A* using Manhattan

A* using Euclidean

A* Algorithm – Heuristic Function



A* using Chebyshev

- All three heuristics found same optimal solution but did not search the same set of nodes
- Blue nodes represent closed set.
- Green represent open set.
- White nodes were not searched.

A* Algorithm - Process

- To actually perform the search we start with two sets
 - List of open nodes to be searched
 - List of closed nodes that have been searched
- Add starting node to open list initially
 - Expand open set to include each neighbor of starting node
 - Put starting node on closed set
 - Calculate the F costs of those open neighbor nodes
- During the next round we chose the neighbor node with the lowest calculated F cost to expand next and repeat the algorithm, until either a solution is found or we exhaust the set of available nodes

A* Algorithm – Full Pseudo Code

```
initialize the open list
initialize the closed list
put the starting node on the open list (you can leave its f at zero)
while the open list is not empty
  find the node with the least f on the open list, call it "q"
  pop q off the open list
  generate q's 8 successors and set their parents to q
  for each successor
           if successor is the goal, stop the search
     successor.g = q.g + distance between successor and q
     successor.h = distance from goal to successor
     successor.f = successor.g + successor.h
     if a node with the same position as successor is in the OPEN list \
       which has a lower f than successor, skip this successor
     if a node with the same position as successor is in the CLOSED list \
       which has a lower f than successor, skip this successor
     otherwise, add the node to the open list
  end
  push q on the closed list
end
```

A* Algorithm – Common Assumptions

- The graph to be searched is usually a grid with uniform cost to move from node to node
- Optimal data structures like priority queue are available
- Certain nodes act as obstacles that cannot be traversed easily if at all
 - Examples: wall, bodies of water



Our go to visual example with walls and relocated end point

A* Algorithm – Running Time

- Under ideal circumstances the amount of nodes searched relative to the graph size is small
- At worst every node must be searched to find solution as well as when the solution doesn't exist



Near worst case performance. Also note solution found is not optimal!!!

A* Algorithm – Parallelizing

• Need to find a method to properly utilize parallel processing environment

 There are lots of variations on the A* algorithm that introduce a parallelizable component

 We are going to look at the hierarchical breakdown method and show that it allows for a parallel prefix style solution

Hierarchical Breakdown A* (HBA)

- Adds a level of abstraction over graph
- Splits graph into clusters
- Algorithm finds entry and exit points between clusters



Our base map before HBA divided into four clusters

- Now use algorithm to find points that connect the clusters
- These points are marked in yellow



- In each cluster use A* to find the distances between each yellow node
- We know the cluster that contains our start and end nodes
- Use local A* to move from start node to HBA paths
- Travel along optimal HBA paths until end node cluster reached
- Use local A* to travel to end



- Path not guaranteed optimal, but within 1%
- Trade off of optimal solution for computation time and smaller memory footprint
- Useful as a preprocessing step that can be reused for multiple path searches on same map



Path taken by base A* using Chebyshev heuristic. Compare to HBA path.



Parallelizing HBA

- We break down the graph into clusters during HBA
- These clusters can be given to our set of PEs
- Compute entry points and the paths between them locally on each PE
- We know cluster location of goal node
- Each cluster calculates how to reach desired goal cluster using HBA paths
- PEs communicate these sub solutions to each other until all nodes know a solution (our parallel prefix step)
- Use MPI to accomplish message passing and internode communication

Experimental Setup

- Multiple map types
 - Empty rooms
 - Spiral
 - Maze-like
 - Open area with large obstacles
- Test on various sizes
 - 512x512
 - 1024x1024
 - 4096x4096
 - 8192x8192

Experimental Setup

- Change cluster sizes for each map
 - 16x16
 - 32x32
 - 64x64
- Vary number of PEs used
 - 1,2,4,8,16,etc

Examples of "Professional" Maps Used





Open Area Map

- Map used as a test sample
- Will run tests on this first including changes to
 - Overall map size
 - # of PEs used



Results – Open Map

)9
3
4
4
3
3
3

Map Size

All results using 32x32 HBA cluster sizing

Results are running time in milliseconds

Experiment includes time to set up HBA clusters and navigate a path spanning the map

Realistically the time spent making HBA clusters can be reused for repeated path calculations resulting in overall dramatically reduced times

Results – Open Map

Yending 12000 10000 8000 4000 4000 **→**8192 **---**4096 -1024 **--**512

Running Time vs #PEs

Results – Open Map



Results – Open Map Speedup

			Map Size		
		512	1024	4096	8192
#PEs	2	1	1	1	1
	4	2.019848708	1.530999255	1.880822819	1.495367095
	8	3.441206661	3.091925397	3.464122177	1.997923925
	16	4.889798234	8.193711135	4.465605843	3.070385947
	32	2.560571986	17.25492618	5.686721431	3.210998699
	64	1.947555269	12.70766991	3.829731544	2.906496515
	128	1.918393947	11.67126264	2.748024959	2.552535595

Results – Open Map Speedup

Speedup vs #PEs



Conclusions

- HBA allows for an easy implementation of the A* search algorithm
- There are many solutions to parallelizing that can be explored though
- Thorough experiments take time to conduct
 - Did not have enough time/resources to generate data for more than one map type
- Running times increase drastically as map size increase
 - More efficient data structures needed for maintaining node lists
 - Message passing large datasets in costly
- Speedups follow a similar trend though
 - For this setup maximum speedup was achieved with using 32 nodes for computation

References

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References

- Visual examples created using PathFinding.js.
 - http://qiao.github.io/PathFinding.js/visual/