

# Assignment 1

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198:520 Intro to Artificial Intelligence  
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1) Code is attached for question 1.

2)

2.1) In order to choose a map size that allows us to experiment and reach valid theoretical conclusions, we had a few criteria that needed to be met:

1) The map size had to be small enough so that we could visually see the maze on our screens.

2) The runtime of each algorithm over 100 iterations would remain around 1 second so that we could test a wide range of  $P$  values.

3) The time it would take a human to solve the maze would be roughly between 10-20 seconds.

Based on these criteria, we found **dim** = 30 sufficient to satisfy our needs. All tests were done with this dimension unless noted otherwise.

2.2) Our implementation of BFS adds unblocked/unvisited cells to a queue, checking that a given cell's neighbors are unblocked/unvisited in the following order: right, down, up, left. Cells in the queue are checked in the order that they are added until the goal cell is reached. Given the enqueue/push ordering of each different algorithm, we can see (in the visual below) that these algorithms work as intended.

Our implementation of DFS adds unblocked/unvisited cells to a stack, checking that a given cell's neighbors are unblocked/unvisited in the following order: up, left, right, down. Cells in the stack are checked with the most recently added cell being checked first until the goal cell is reached. As a result, it makes sense that the final path favors going down and going right, in that order.

Our implementation of A\*-Manhattan adds unblocked/unvisited cells to a queue, ordering the queue based on the lowest estimated distance from the start to the goal using the sum of the distance traveled to a cell and the Manhattan distance from the cell to the goal

and checking that a given cell's neighbors are unblocked/unvisited in the following order: right, down, up, left. As a result, it makes sense that the final path limits the Manhattan distance from the start cell to the goal cell.

Our implementation of A\*-Euclidean adds unblocked/unvisited cells to a queue, ordering the queue based on the lowest estimated distance from the start to the goal using the sum of the distance traveled to a cell and the Euclidean distance from the cell to the goal and checking that a given cell's neighbors are unblocked/unvisited in the following order: right, down, up, left. As a result, it makes sense that the final path limits the Euclidean distance from the start cell to the goal cell.

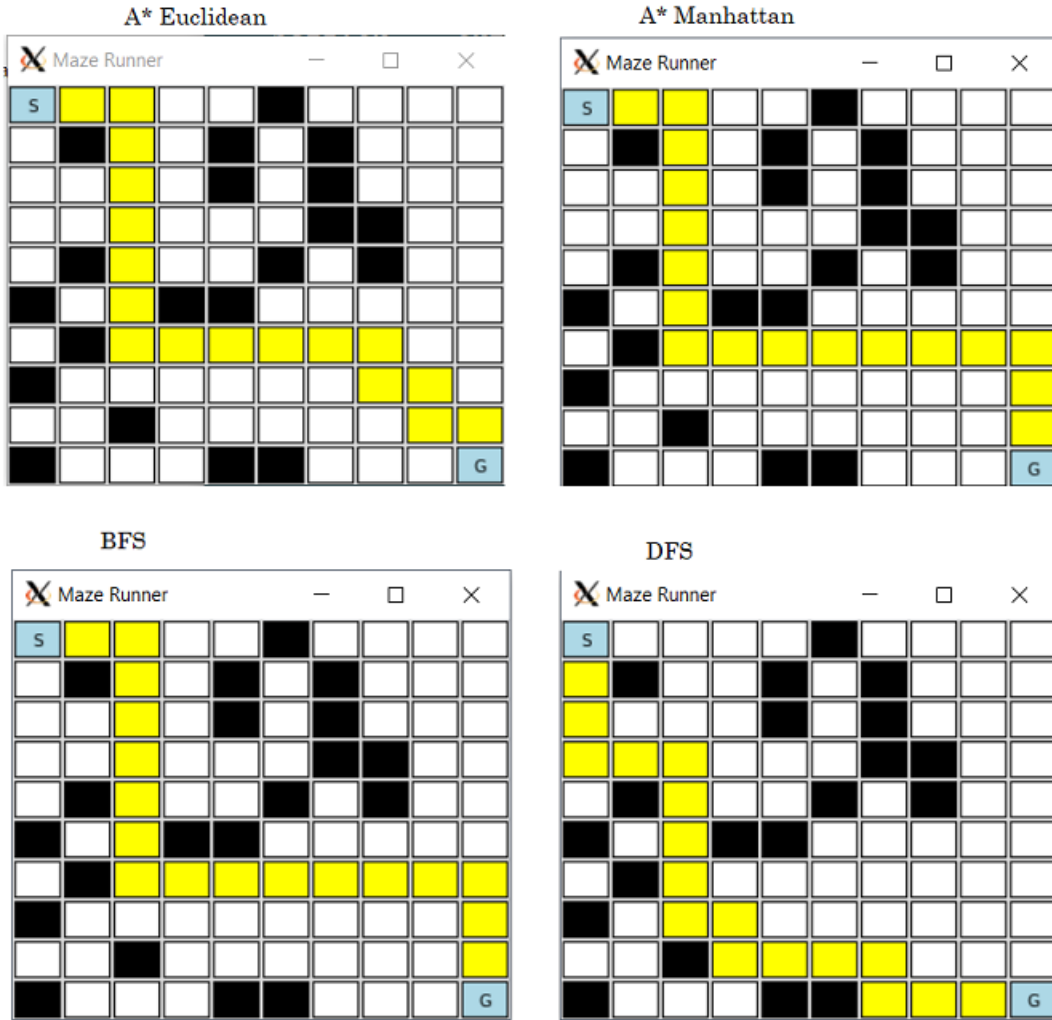


Figure 1: Solution Paths

**2.3)** Given  $\text{dim} = 30$ , Figure 2 depicts how maze-solvability depends on the value of  $p$ .

Methodology to generate graph for  $p = 0.0, 0.1, 0.2, \dots, 0.9$  : run DFS on a maze with  $\text{dim}$

30. Repeat this for a total of 100 tests and keep track of how many were solvable. Then that ratio will give the estimate for the probability that a maze of the corresponding dim and  $p$  will be solvable.

The best algorithm to use here is DFS because it runs the fastest. We don't need optimality if all we need to know is if the maze is solvable or not, and DFS is guaranteed to find a path to the goal if one exists. We take 30 to be a good sample size that is representative of the true ratio of solvability for a given  $p$ .

We can see that the maze-solvability drastically decreases as the value of  $p$  increases. The more interesting question is, what is the threshold value,  $p_0$ , where for all  $p < p_0$ , most mazes are solvable (here we define most as any value greater than 50%), and for all  $p > p_0$ , most mazes are not solvable. Based on the tests and the results depicted in this graph, that threshold value is  $p_0 = 0.3$ .

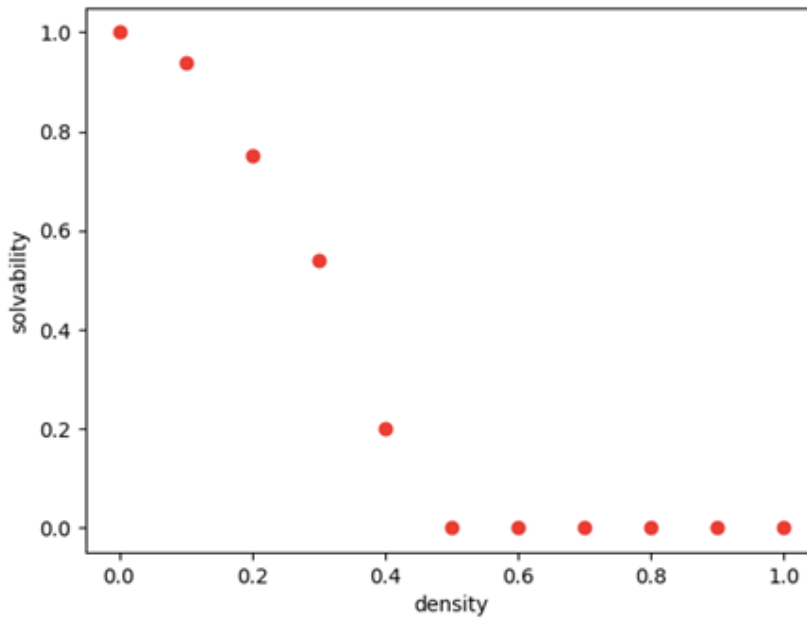


Figure 2: Solvability vs. Density

- 2.4) For each  $p$  in  $[0, p_0]$ , we generated and solved 30 mazes, and found the average shortest path (see Figure 3). We chose a step size of 0.2. The BFS and A\* algorithms are the most useful for this problem, because they return the shortest path by default.

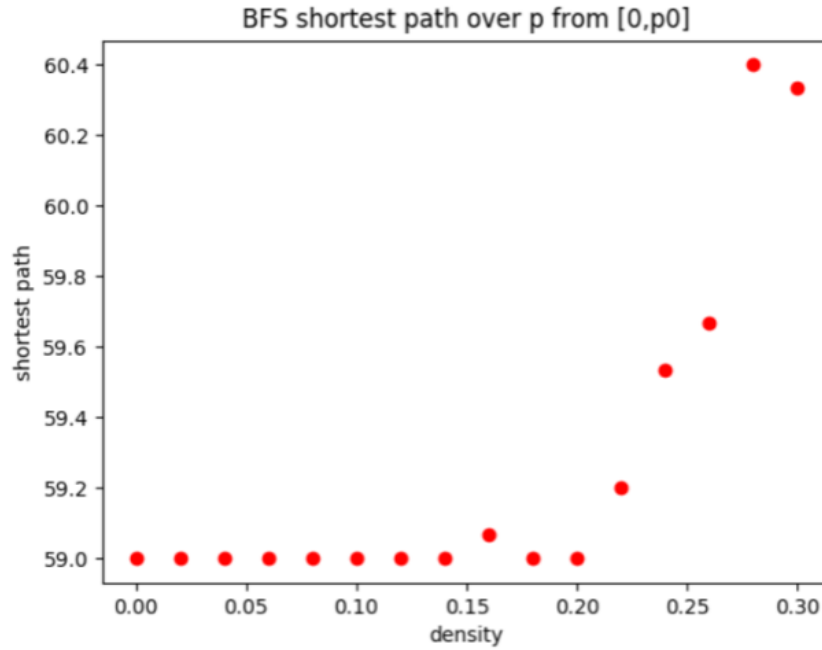


Figure 3: Expected Shortest Path vs. Density

- 2.5) The Euclidean distance heuristic helps produce a more direct path to the goal, however it is more computationally heavy. The Manhattan distance heuristic leads to more stretches of straight line paths within the overall final path, or in other words - less frequent changes of direction (less zig-zagging), but the path from the start to the goal is not necessarily direct and almost linear, as the Euclidean heuristic would have provided us.

Computing A\* with the Euclidean distance heuristic takes much longer than computing with the Manhattan distance because it typically requires visiting more nodes. This is because when we estimate  $h(n)$  using Euclidean distance, that results in lower values for  $h(n)$ , at some nodes, than they would have been with the Manhattan distance heuristic, therefore, we are more likely to visit those nodes.

Figure 4 shows how the two algorithms behave in comparison to each other, in terms of nodes expanded and runtime, with respect to  $p$ . In the left graph of Figure 4 the average runtime of A\* Euclidean is vastly greater than the average runtime of A\* Manhattan, for  $p$  values that yield a high percentage of solvable mazes.

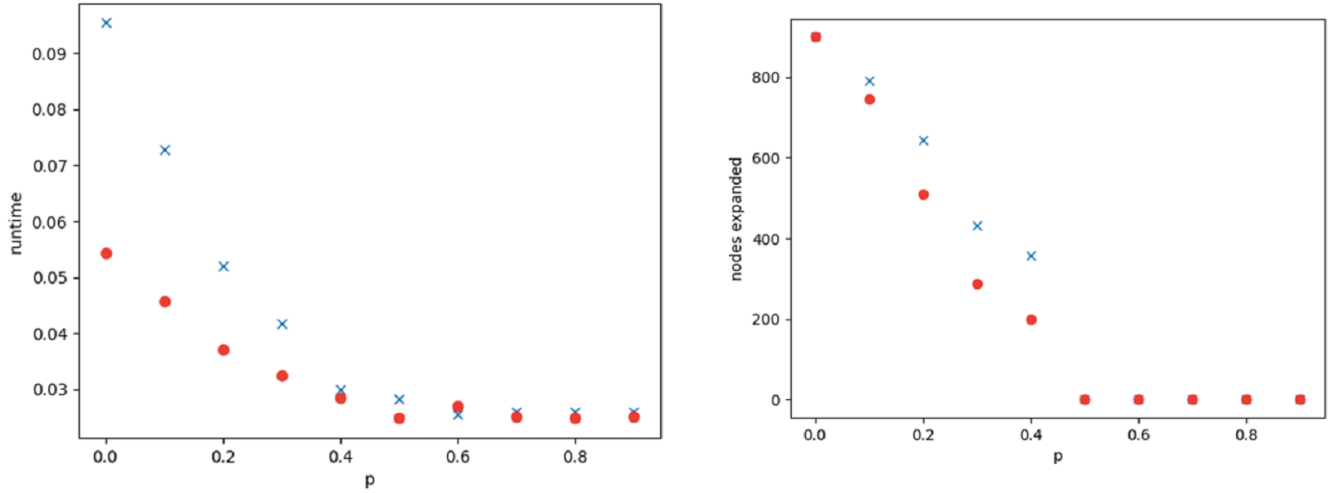


Figure 4: Manhattan vs. Euclidean (Red dot: Manhattan, X: Euclidean). Note: runtime measured in seconds

This is supported by the results shown in the right graph in Figure 4, where we can see that A\* Euclidean expands more nodes on average, than A\* Manhattan does, (for  $p$  values that yield a high percentage of solvable mazes) except for when  $p = 0.0$  because both algorithms then expand every node in the maze.

So, depending on the problem we want to solve, there will be trade offs for using either heuristic. The Euclidean distance is useful to generate more direct paths, but at the cost of increase computation time. The Manhattan distance is useful if you want quicker computation, but will give you paths that are suited for a grid-like area and not necessarily a straight-shot from start to goal with zero changes in direction.

- 2.6)** We generate a graph for average path length at various  $P$  to show that BFS is better than DFS at finding the shortest path. For the left graph in Figure 5, we see that when  $P$  is 0, BFS and DFS both produce solutions of equal length. However as  $P$  increases, the average solution lengths produced by DFS get longer, while the lengths produced by BFS remain stable (around 60). We also generated a scatter plot (the right graph of Figure 5) to represent the difference in length of individual maze solutions. The y-axis represents the length difference between DFS and BFS paths. We can see that there are no instances where the difference is less than 0 (we represent positive differences with red dots, and negative ones with black dots), thus there are no instances where DFS finds a better path than BFS.

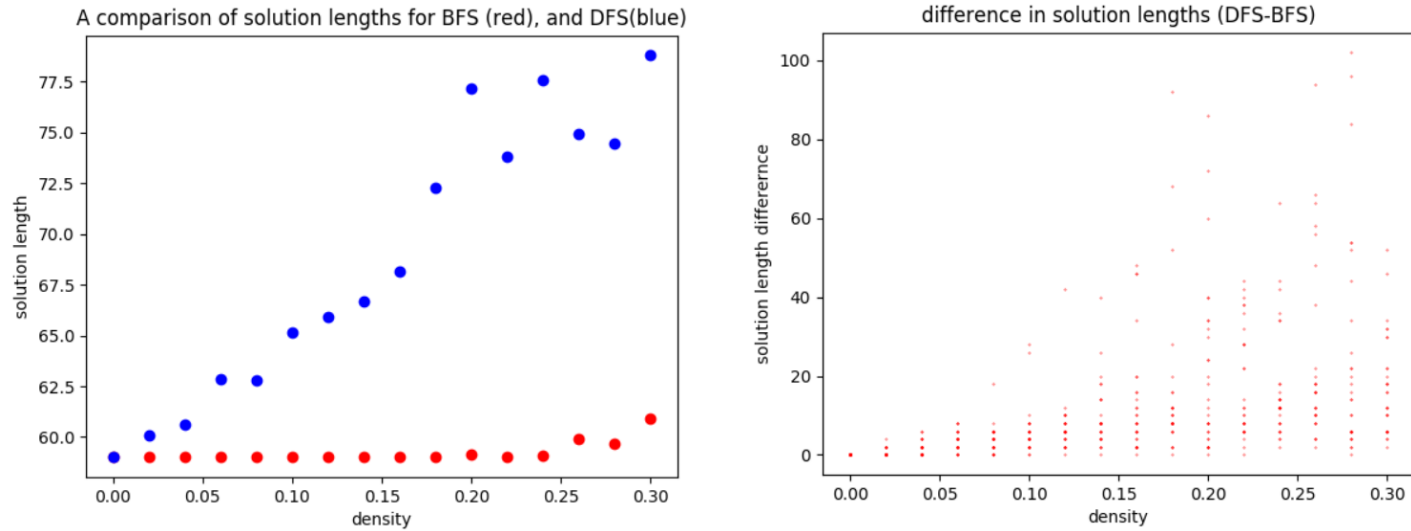


Figure 5: BFS vs. DFS

- 2.7)** All four algorithms behave as expected. BFS uses a queue to find a path from the start cell to the goal cell, in addition to being optimal. DFS uses a stack to find a path from the start cell to the goal cell, and as a result of the nature of the algorithm and its implementation, is not optimal. A\*-Manhattan and A\*-Euclidean are modified versions of BFS, ordering the queue such that the lowest estimated distance from the start cell to the goal cell through a given cell is prioritized within the queue. However, the differing heuristics result in differing final paths for the two algorithms.
- 2.8)** DFS performance can be improved by modifying the order in which neighboring cells are added to the fringe. For example, neighboring cells can be added to the stack such that the closest cell to the goal cell is the most recently added cell, so that each set of additions to the fringe prioritizes getting closer to the goal cell. DFS is not optimal, and as a result can expand more nodes than is necessary, so it is important that any modifications to the order of adding neighboring cells to the fringe intend to limit the number of nodes expanded and return a shorter path to the goal cell.
- 2.9)** The threshold probability  $p_0$  varies as dim changes. At lower values of dim ( $\text{dim} < 30$ ), the majority of mazes become unsolvable at lower values of  $p$ , typically near 0.24 - 0.27. Conversely, as you increase the value of dim ( $\text{dim} > 30$ ), the threshold value,  $p_0$ , where most mazes become unsolvable, converges at 0.3.
- 3)** We decided to use beam search to perform local search. We represented the maze as a 2D list of Cell objects, which contain 1s and 0s. Each maze that we generate is solved and assigned a difficulty level, and our goal is to find the maze with the highest difficulty for its respective search and metric. Our implementation of beam search involved generating 8 initial mazes. Out of each of those mazes, we run it through an edit-maze function, which randomizes the 1s and 0s in each maze with a designated probability (we set it at 0.1). This process generates 2 child

mazes, which we calculate the difficulty of. After we have a list of 16 mazes, we select 8 of those mazes to move onto the next level, until we terminate at a fixed number of levels. In an attempt to maximize genetic diversity, we selected  $0.75 \times 8$  of the best mazes and randomly selected  $0.25 \times 8$  of the other mazes to move onto the next level (generation).

In order to code beam search to generate hard mazes, we had to account for many parameters: beam length, the edit probability, the number of child mazes generated by each maze, the percentage of best/worst mazes to propagate, how many generations of mazes to create, the initial starting probability of the generated mazes. We had to decide how to select mazes for the next generation to optimize diversity and difficulty

A problem that we faced was the issue of genetic diversity. This was tested by running our algorithm using BFS to find maximal shortest path. Intuitively, this should generate a maze that looks like Figure 6:

|    |    |    |    |    |    |    |    |    |
|----|----|----|----|----|----|----|----|----|
| 0  | 1  | 2  | 3  | 4  | 5  | 6  | 7  | 8  |
| 9  | 10 | 11 | 12 | 13 | 14 | 15 | 16 | 17 |
| 18 | 19 | 20 | 21 | 22 | 23 | 24 | 25 | 26 |
| 27 | 28 | 29 | 30 | 31 | 32 | 33 | 34 | 35 |
| 36 | 37 | 38 | 39 | 40 | 41 | 42 | 43 | 44 |
| 45 | 46 | 47 | 48 | 49 | 50 | 51 | 52 | 53 |
| 54 | 55 | 56 | 57 | 58 | 59 | 60 | 61 | 62 |
| 63 | 64 | 65 | 66 | 67 | 68 | 69 | 70 | 71 |
| 72 | 73 | 74 | 75 | 76 | 77 | 78 | 79 | 80 |

Figure 6: Optimal Difficulty

We got close, but we were never able to generate a maze that looked as optimal as above, even after the maze difficulties plateaued (which is our termination condition). This implies that our maze generation algorithm leads us to become stuck at a local maxima, despite our efforts to maintain genetic diversity. We might be able to optimize this by reducing the percentage of best performers that move onto the next generation, or by increasing the edit probability parameter. The termination condition we applied was to print the current maximum difficulty after each generation of mazes, and eyeball it until the difficulty stopped increasing. The advantage of this

heuristic is that our intuitions are often correct about when to stop running the algorithm. The disadvantage is that this would be inconvenient for running a large number of trials. Ideally we could experimentally determine a function for when to stop, that varies based on the dimension of the maze. Another disadvantage is that we might be at a local maxima of difficulty, and terminating early might lead to never finding a higher difficulty where one exists.

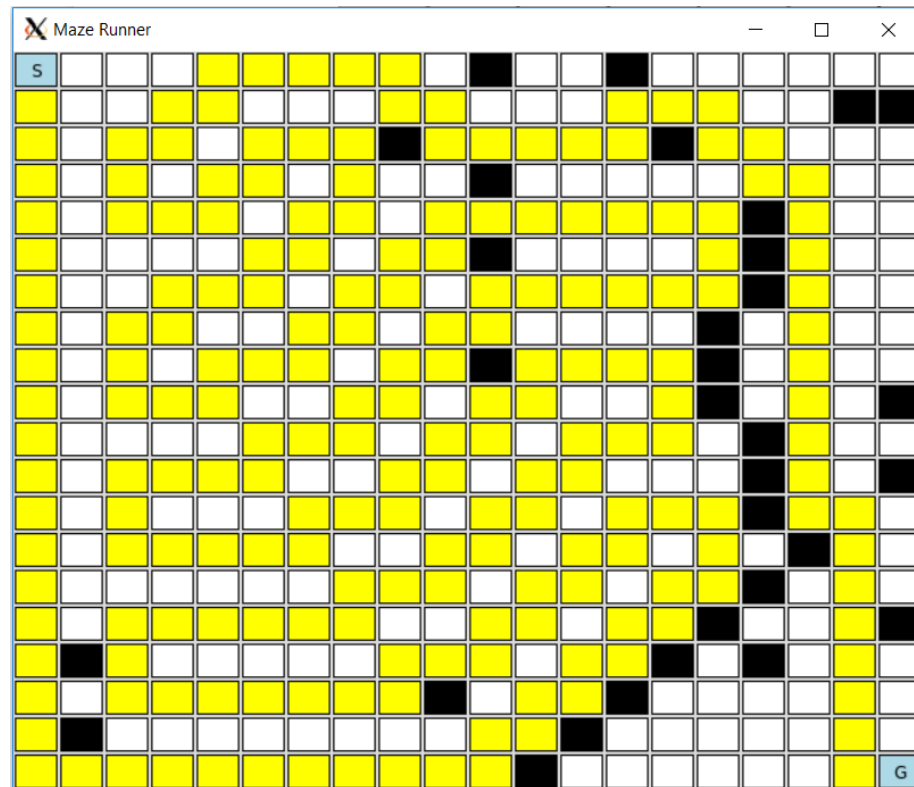


Figure 7: DFS with shortest maximal path

DFS with maximum shortest path:  
 Solution length: 199  
 Max Fringe: 136  
 Max Nodes: 337

(Figure 7) Initially this did not agree with our intuition, because one would expect a difficult maze to have more obstructions. The structure of this maze is a representation of the limitations of DFS more than the shortest existing possible path through the maze (which can be found using BFS). The back and forth snaking through the empty sections of the maze reflects the order in which we put nodes into the fringe.



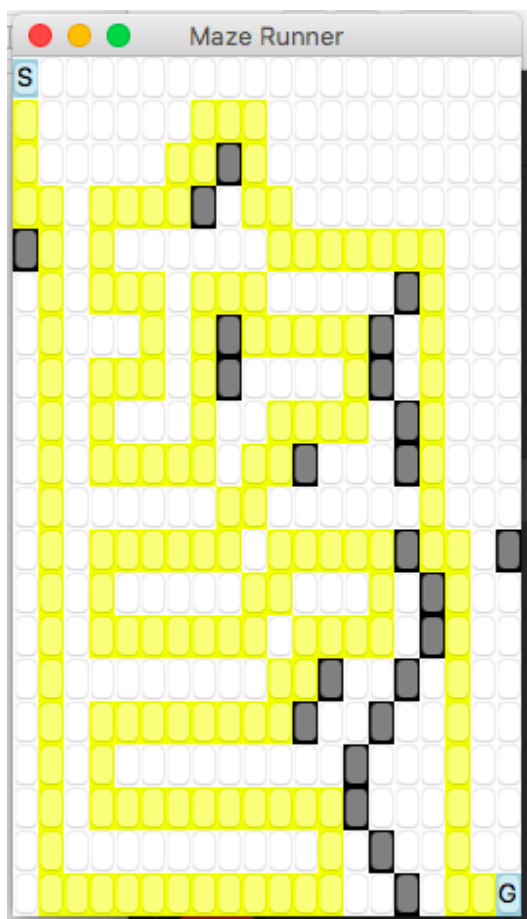


Figure 8: DFS with maximal fringe size

DFS with maximal fringe size:

Solution Length: 139

Max Fringe: 144

Max Nodes:  $\approx 350$

(Figure 8) The hard maze generated by the paired metric "DFS with max-fringe size" agreed with our intuition for the following reasons. Following similar logic to the A\* Manhattan with max-fringe metric, a long zig-zagging path that avoids the edges of the maze is what we expect to see and is exactly what we got. This is conducive to "seeing," and consequently, incorporating more neighboring nodes into the fringe. Also, this DFS path looks similar to the DFS path generated for the maze that was created with the "DFS with maximal shortest path" paired metric, which makes sense, because fringe size is correlated positively with path length. The key differences in the mazes generated using the two different metrics for "hard" is that the max fringe size metric produced a maze with fewer obstructions. This makes sense, because that allows the fringe to fill up as the algorithm will find that more of the neighboring nodes are eligible since they are not blocked.

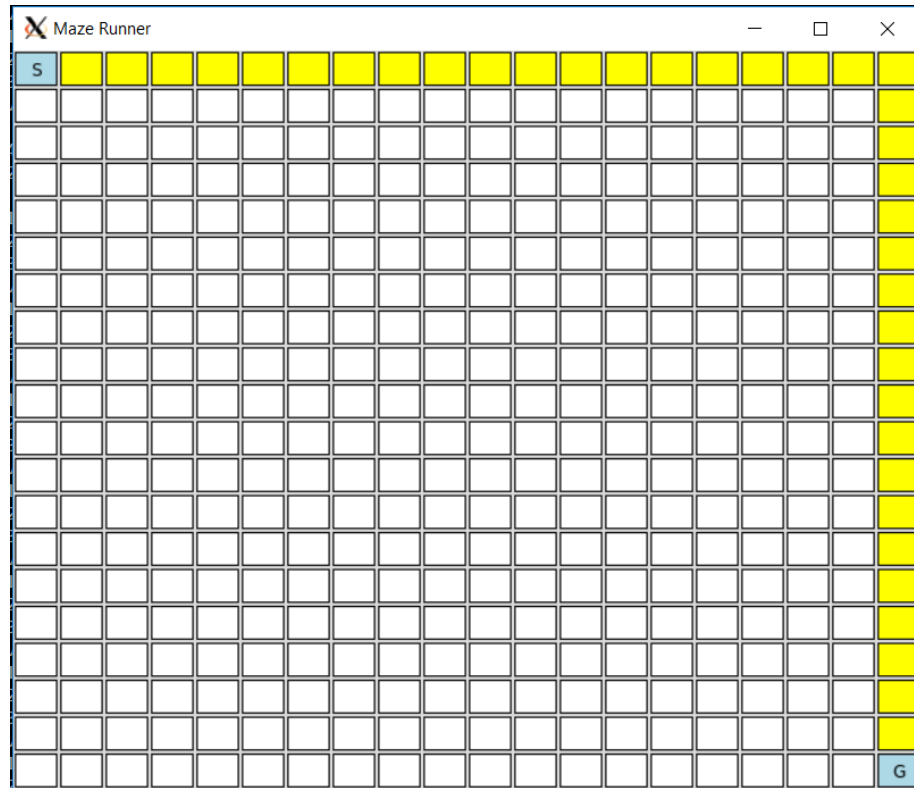


Figure 9: A\* Manhattan max nodes expanded

A\* Manhattan max nodes expanded:

Solution Length: 39

Max Fringe: 21

Max Nodes: 400

(Figure 9) This agrees with our intuition. Because there are no blockades, every possible position has the same  $g(x) + h(x)$  value, where  $g(x)$  is length so far and  $h(x)$  is the heuristic. This means that our priority queue devolves into a normal queue, which leads to a normal BFS being performed. BFS will expand to every possible node in between the start and the goal, so it makes sense that 400 nodes are expanded for a 20 x 20 maze.

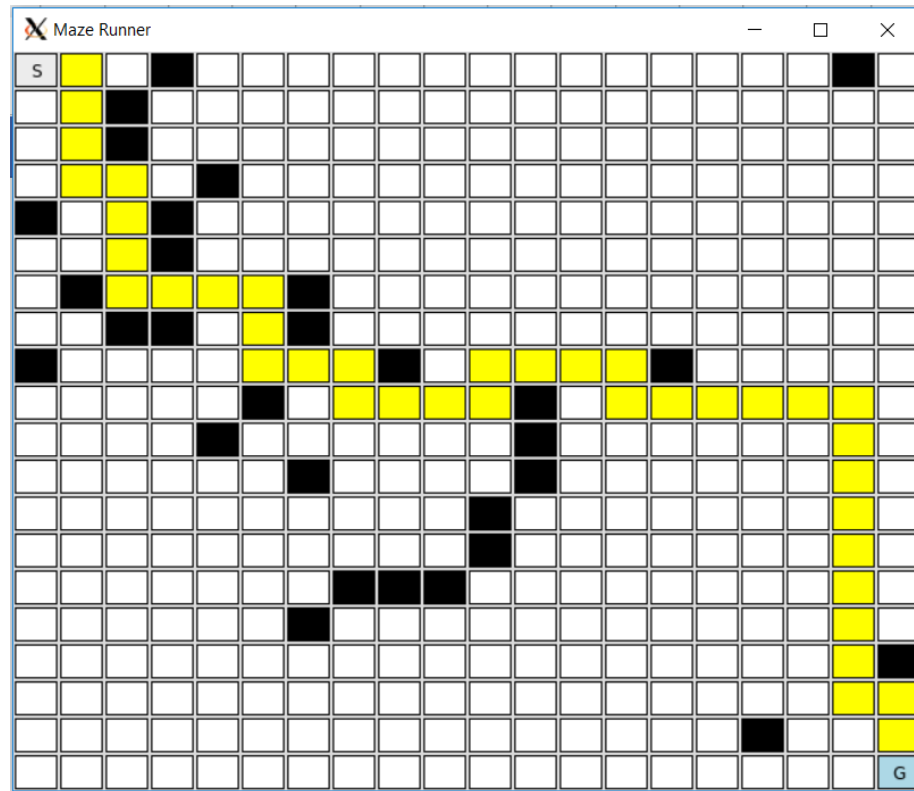


Figure 10: A\* Manhattan max fringe size

A\* Manhattan max fringe size:

Path len: 41

Max Fringe Len: 97

Max Nodes Expanded: 354

(Figure 10) This solution initially did not match our intuition because we were expecting a maze that was void of any obstructions, since we thought that would cause the maximum number of nodes to fill up in the fringe at one time. However, after some thought this result seems to make sense because if the algorithm's path went only on the edges of the maze, it wouldn't necessarily "see," and hence, wouldn't add as many nodes to the fringe because it ran along the border of the maze. If you have a longer, snaking path within the maze, maybe one that even goes away from the goal for a few steps (i.e. up or left a couple times, especially since the order we add nodes to our fringe is in a clockwise fashion - that is, add right neighbor first, then bottom neighbor, then left neighbor, then top neighbor), while having neighbors on both sides (unlike along the edges of the maze), you'll pick up more nodes for the fringe as you explore. Additionally, because this is an A\* algorithm, it was not going to explore many of the nodes that are off of the optimal path anyway, that's why we see that the obstructions are mainly near the top left of the maze, so we are close to one of the largest fringe sizes possible with A\* manhattan. Also, having obstructions helped in two ways: one, it kept the algorithm away from just crawling along the edges of the maze toward the goal; two, it caused the algorithm to snake around to more nodes which allows

it to pick up more neighbors that it would if it just went to the right and then down without changing direction ever. This also extends the size of the final path, which will also cause the fringe to grow.

4) Intuitively, the notion of a thinning A\* seems like a viable efficient approach to searching. We found, however, that this is not the case. We sought to find the overall cost of using a thinning A\*, using A\* Euclidean as our base search algorithm, when compared to using A\* Euclidean without the thinning process. We define cost as the total amount of nodes the search algorithm must expand. The thinning A\* approach can be described as a sort of optimal directed DFS search; we solve a simpler maze, use the solution path as our heuristic for solving the original maze, and take detours if blocked cells are encountered along the path. This process by itself is faster than running any of the previously used search algorithms, yet when combined with the cost of obtaining the heuristic, this technique becomes less efficient as can be seen in Figure 11.

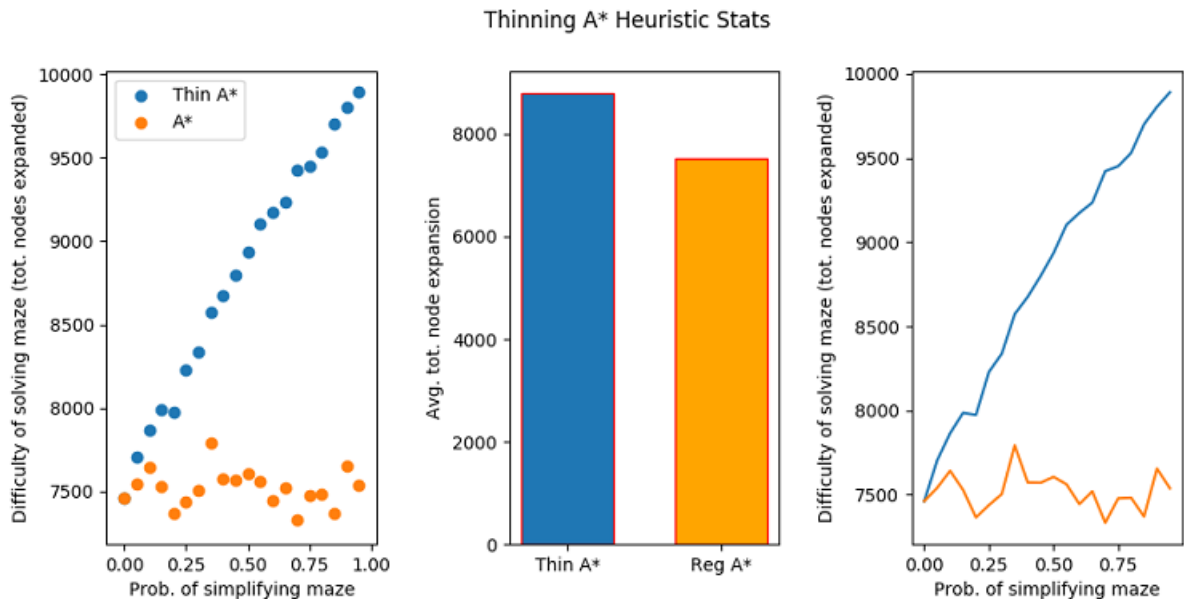


Figure 11: Thinning A\* Heuristic

When considering the cost of computing the thinning A\* heuristic, we quickly run into a problem; the overall cost of computing this heuristic get significantly larger, in terms of overall nodes expanded, as the probability ( $q$ ) of simplifying the maze gets larger. That is, the easier the maze we solve, the more work the thinning algorithm has to do. It seems that having blocked cells in a given maze reduces the amount of work needed to search through that maze because these blocked cells constraint the search space - they eliminate paths from the search tree. A thinning algorithm can still be valuable, however, if the thinned maze is  $\approx 5\%$  easier than the original. It is worth noting that we did not evaluate all the possible way this heuristic can be used or calculated and thus there might be other ways to exploit this idea. For example, if we know that the shortest path of the thinned maze is 10 cells, then we might be be able to use this as a

lower limit when solving the original maze by DFS limiting the search depth to 10 or so. Yet, we would still run into the problem of computing the heuristic.

The heuristic by itself is useful, yet the cost of calculating it makes it comparable, and usually less favorable, to the other A\* algorithms for this specific problem.

5) In the previous examples our main goal was to find an optimal path, shortest in most cases, from start to finish. In this problem, however, there is a question of safety versus optimality. We are given that the probability of a cell catching on fire is  $P(A) = 1 - (\frac{1}{2})^k$  where  $k$  is the number of neighbor cells on fire.

For clarity, we assume that neighbors only count as adjacent cells and not diagonals. The main idea behind the proposed algorithm (shown below) is creating a safest/optimal using a safety heuristic - the robot, or player, will move to any neighboring cell with a probability of  $P(A^c) = 1 - (1 - (\frac{1}{2})^k)$ . Essentially, this is a greedy best first search where the safest neighbor will be prioritized. We propose the following algorithm:

**Input:** A maze (two-dimensional array)  $M$  and starting vertex  $v$  (this is the starting cell) with blocked, fiery, or open cells as described by the problem .

**Output:** A valid path through the maze - a list  $L$  of coordinates.

```

1) Let fringe be a queue
2) enqueue  $v$  onto fringe
3) while fringe is not empty
    set node = fringe.dequeue()
    if node is goal state
        backtrack path to  $v$  through parent nodes and return list of cell coordinates  $L$ .
    else if node has not been visited
        label node as visited and record parent
        for all child nodes (neighboring cells) do
            discard any node that is blocked or on fire
            evaluate safety heuristic  $P(A^c)$  for each open node
            enqueue nodes from safest to least safe (highest to lowest probability)*
        else (node has been visited before or is on fire/blocked)
            discard node and continue
4) if no path is found : return none
note:* if nodes have equal safety heuristic value, enqueue in any order (clockwise, left to right, counterclockwise...).

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### Work Distribution

- Q1 - Split evenly between all group members
- Q2 - Each group member had roughly 1-2 subquestions
- Q3 - Split evenly between Andrea and Karthik
- Q4 - Split evenly between Farhat and Matan
- Q5 - Matan

Overall each member of the group roughly contributed equally to the success of the project.