

# CS 1240 Programming Assignment 1: Random MSTs

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Spring 2026

## 1 Overview

We implemented minimum spanning tree (MST) algorithms and ran experiments on the random graph models described in the assignment. For each  $n$ , we generated multiple independent random instances and recorded the average MST weight. We then fit a simple function  $f(n)$  describing how the expected MST weight grows with  $n$ .

Program interface: `./randmst 0 numpoints numtrials dimension`

Dimensions:

- Dimension 0: complete graph with i.i.d. weights in  $[0, 1]$
- Dimension 1: hypercube graph with i.i.d. weights in  $[0, 1]$
- Dimensions 2,3,4: complete graphs on random points in  $[0, 1]^d$  with Euclidean weights

## 2 Algorithms and Implementation

### Deterministic Edge Weights (Dimensions 0 and 1)

For dimensions 0 and 1, we do not store edge weights explicitly. Instead, we compute weights on demand using a deterministic hash of  $(\text{seed}, u, v)$ . For a fixed seed, this defines a fixed weighted graph, ensuring consistency within a trial while avoiding storage of all edges.

### Dimension 0: Dense Complete Graph

We implemented the standard array-based version of Prim's algorithm.

At each iteration:

1. We scan all vertices not yet in the tree to find the smallest connecting edge.
2. We update the best known connection for every remaining vertex.

Runtime per trial:  $O(n^2)$  Memory usage:  $O(n)$

### Dimension 1: Hypercube Graph

Each vertex  $v$  is connected to vertices of the form  $v \pm 2^i$  for powers of two  $2^i < n$ . Thus each vertex has at most  $\log_2 n$  neighbors, and the graph contains  $\Theta(n \log n)$  edges.

We implemented Prim's algorithm using a custom binary min-heap with decrease-key support. Each vertex appears exactly once in the heap. When a better connecting edge is found, we update its key and restore heap order.

Since each vertex has  $O(\log n)$  neighbors and each decrease-key costs  $O(\log n)$  time, the total runtime per trial is:

$$O(n(\log n)^2).$$

### Dimensions 2–4: Geometric Complete Graphs

We generate  $n$  random points in  $[0, 1]^d$ . Distances are computed on demand using the Euclidean metric.

We use the dense array-based version of Prim’s algorithm, resulting in:

$$O(n^2)$$

runtime per trial.

## 3 Experimental Results

### Dimension 0

$n$	trials	average MST weight
128	5	1.13502
256	5	1.17091
512	5	1.22843
1024	5	1.23670
2048	5	1.21850
4096	5	1.19779
8192	5	1.19745
16384	5	1.19904
32768	5	1.20348

**Fit:** The average converges to a constant near

$$f(n) \approx 1.20.$$

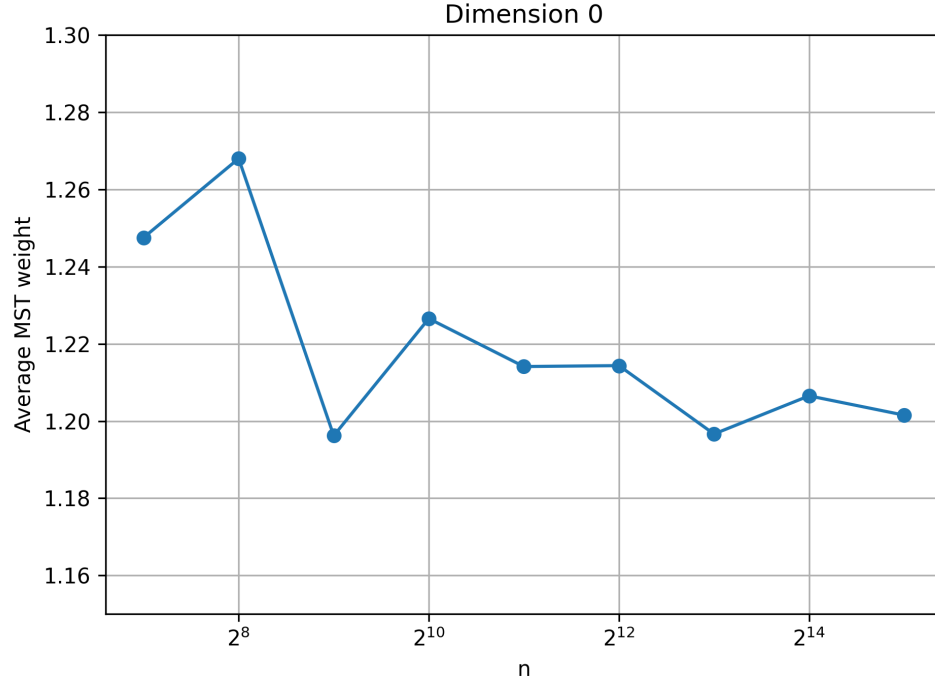


Figure 1: Average MST weight vs.  $n$  for Dimension 0 (complete graph with i.i.d. weights).

### Dimension 1

$n$	trials	average MST weight
128	5	12.2081
256	5	21.3285
512	5	37.8052
1024	5	66.3007
2048	5	121.461
4096	5	219.791
8192	5	404.408
16384	5	742.212
32768	5	1376.88
65536	5	2575.18
131072	5	4842.08
262144	5	9102.08

**Fit:**

$$\frac{9102.08}{262144} \approx 0.0347, \quad f(n) \approx 0.035n.$$

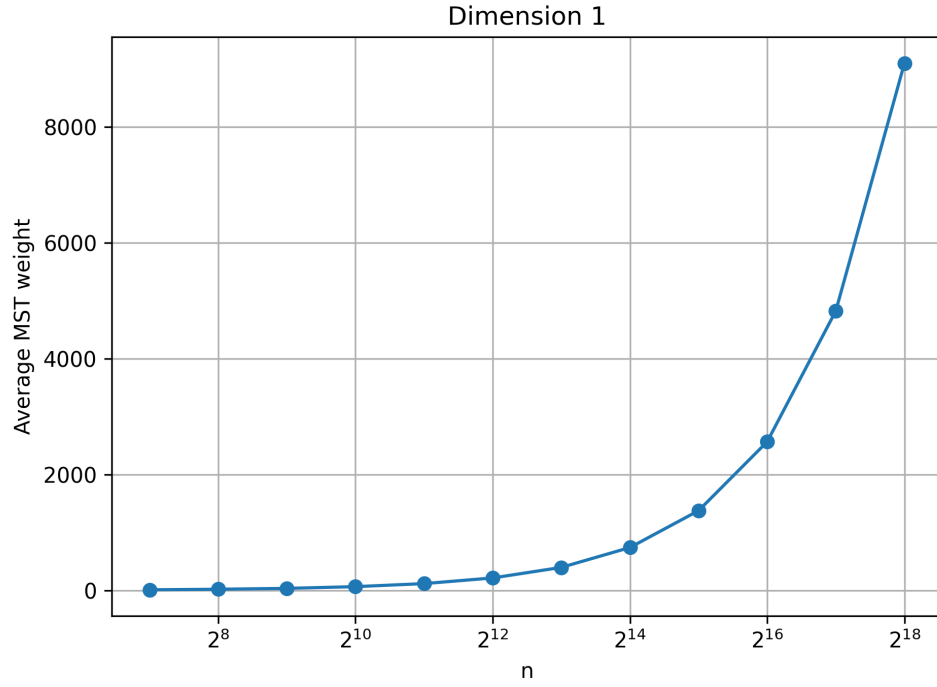


Figure 2: Average MST weight vs.  $n$  for Dimension 1 (hypercube graph).

## Dimension 2

$n$	trials	average MST weight
128	5	7.78429
256	5	10.7068
512	5	14.9969
1024	5	21.1132
2048	5	29.6338
4096	5	41.6704

**Fit:**

$$f(n) \approx 0.66\sqrt{n}.$$

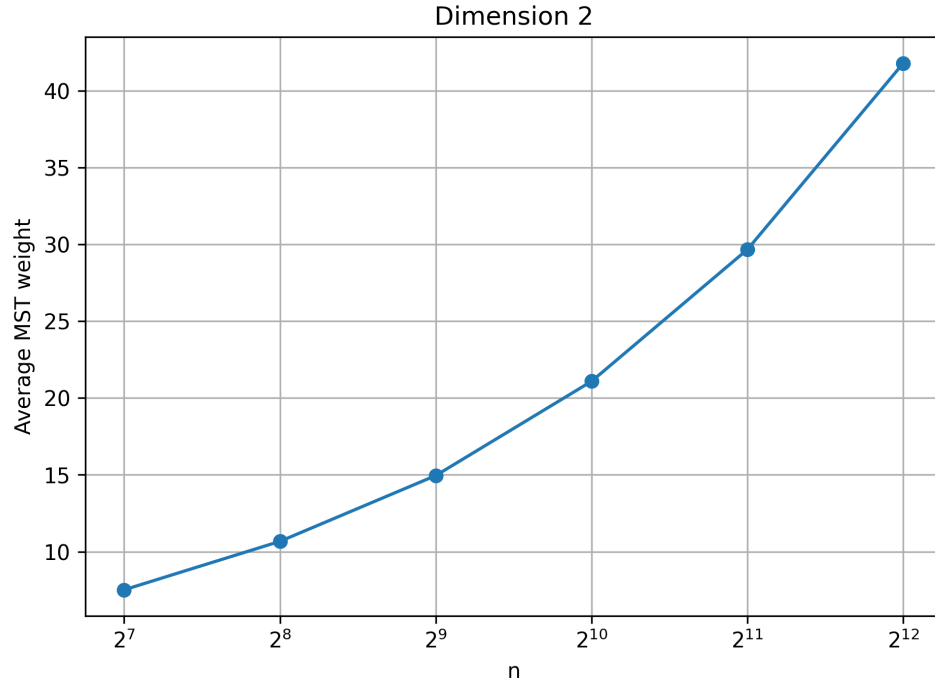


Figure 3: Average MST weight vs.  $n$  for Dimension 2 (random geometric graph).

### Dimension 3

$n$	trials	average MST weight
128	5	17.6731
256	5	28.1207
512	5	43.0517
1024	5	67.8171
2048	5	107.185
4096	5	169.016

**Fit:**

$$f(n) \approx 0.68n^{2/3}.$$

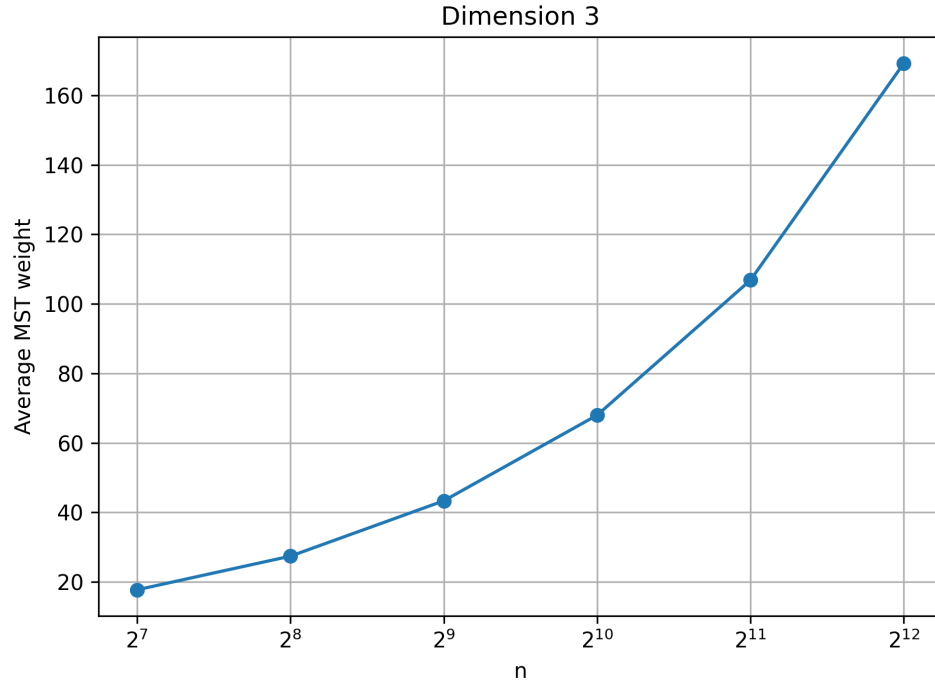


Figure 4: Average MST weight vs.  $n$  for Dimension 3 (random geometric graph).

#### Dimension 4

$n$	trials	average MST weight
128	5	28.3054
256	5	46.6302
512	5	78.2834
1024	5	129.474
2048	5	216.211
4096	5	361.996

**Fit:**

$$f(n) \approx 0.73n^{3/4}.$$

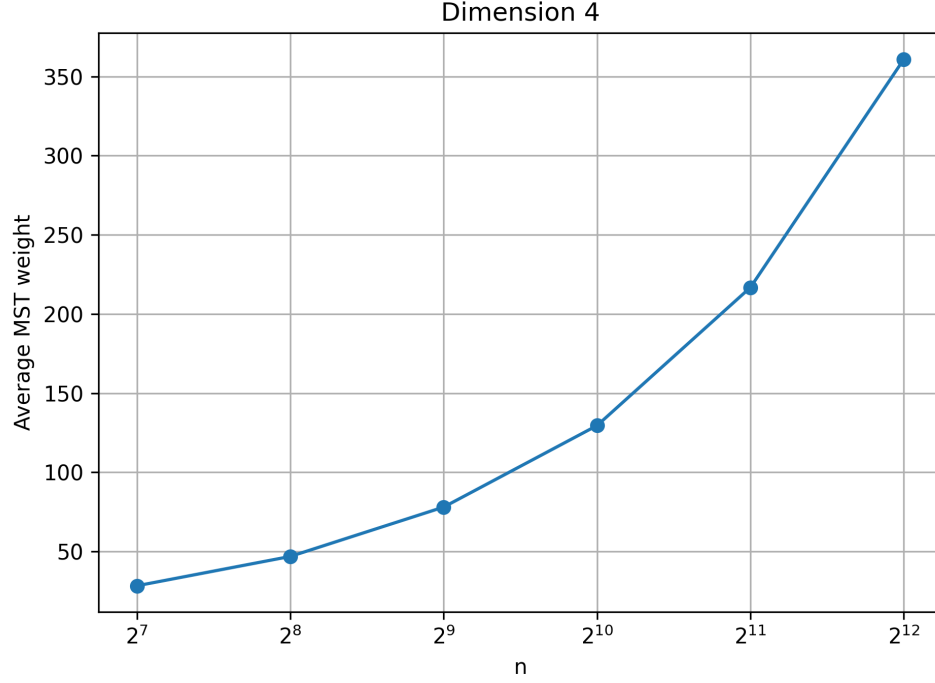


Figure 5: Average MST weight vs.  $n$  for Dimension 4 (random geometric graph).

## 4 Discussion

### Runtime and Scalability

The running time of the implementation depends fundamentally on the structure of the underlying graph.

**Dense graphs (dimensions 0, 2–4).** For dimensions 0, 2, 3, and 4, the graph is complete. A complete graph on  $n$  vertices contains

$$\frac{n(n-1)}{2} = \Theta(n^2)$$

edges. Even though we do not store these edges explicitly, the dense structure forces us to consider  $O(n)$  potential connections at each step of Prim’s algorithm.

In our array-based implementation of Prim’s algorithm, each iteration consists of two operations:

1. Scanning all vertices not yet in the tree to find the minimum key value. This requires  $O(n)$  time.
2. Updating the best-known connection value for every remaining vertex. This also requires  $O(n)$  time.

Since these operations are performed for each of the  $n$  vertices added to the tree, the total runtime per trial is

$$O(n^2).$$

This quadratic growth is unavoidable for dense graphs in this implementation model, because each vertex potentially connects to all others.

For geometric graphs (dimensions 2–4), each relaxation requires computing a Euclidean distance. While this increases the constant factor (due to arithmetic operations and a square root), it does not change the asymptotic complexity.

Memory usage in these cases remains  $O(n)$ , as we store only:

- A boolean array indicating membership in the tree,
- An array of best-known edge weights,
- For geometric graphs, the list of  $n$  points.

**Hypercube graph (dimension 1).** In contrast, the hypercube graph is sparse. A vertex  $v$  is connected only to vertices of the form  $v \pm 2^i$  for powers of two  $2^i < n$ . The number of such powers is at most  $\lfloor \log_2 n \rfloor + 1$ , so each vertex has  $O(\log n)$  neighbors.

Thus, the total number of edges is

$$m = \Theta(n \log n),$$

which is asymptotically much smaller than  $n^2$ .

For this case, we implemented Prim’s algorithm using a custom binary min-heap with decrease-key support. Each vertex appears exactly once in the heap. When a better connecting edge is discovered, we perform a decrease-key operation, restoring heap order in  $O(\log n)$  time.

Each vertex has  $O(\log n)$  neighbors, so the total number of relaxation attempts is  $O(n \log n)$ . Since each successful relaxation requires a decrease-key operation costing  $O(\log n)$  time, the overall runtime per trial is

$$O(n(\log n)^2).$$

This asymptotic improvement over  $O(n^2)$  is what allows the algorithm to scale to  $n = 262,144$ .

## Correctness

All minimum spanning trees were computed using Prim’s algorithm. We now justify its correctness rigorously.

Prim’s algorithm maintains a set  $S \subseteq V$  of vertices already included in the growing tree. At each iteration, it selects the minimum-weight edge that crosses the cut

$$(S, V \setminus S).$$

The correctness of this choice follows from the *cut property*:

**Cut Property.** Let  $(S, V \setminus S)$  be any cut of a weighted graph. If  $e$  is a minimum-weight edge crossing that cut, then there exists a minimum spanning tree that contains  $e$ .

At each step of Prim’s algorithm, the selected edge is the minimum-weight edge crossing the current cut. By the cut property, that edge is safe to add — meaning that it can be extended to a full minimum spanning tree.

Since the algorithm adds exactly  $n - 1$  edges and never forms a cycle, it constructs a spanning tree. Because every added edge is safe, the resulting tree is a minimum spanning tree.

**Deterministic weight generation.** For dimensions 0 and 1, edge weights are not stored but computed via a deterministic hash function of  $(\text{seed}, u, v)$ . For a fixed seed, this defines a fixed weighted graph. Therefore, within each trial, Prim’s algorithm operates on a well-defined weighted graph and produces its exact MST.

Different trials correspond to different seeds and therefore independent random graphs.

## Interpretation of Growth Rates

The experiments reveal distinct asymptotic behaviors that depend on graph structure.

**Dimension 0 (complete graph with i.i.d. weights).** In a complete graph with independent weights uniformly distributed in  $[0, 1]$ , the minimum edge incident to each vertex becomes increasingly small as  $n$  grows. Classical results in probabilistic combinatorics show that the expected MST weight converges to a constant (specifically,  $\zeta(3) \approx 1.202$  for uniform  $[0, 1]$  weights).

Our experiments show convergence toward approximately 1.20, consistent with this theoretical prediction.

**Dimension 1 (hypercube graph).** The hypercube graph is sparse, with only  $O(\log n)$  neighbors per vertex. Unlike the complete graph, increasing  $n$  does not dramatically increase the number of candidate edges available to reduce edge weights. As a result, the average edge weight in the MST does not shrink rapidly with  $n$ .

Since the MST contains  $n - 1$  edges and the typical edge weight remains bounded away from zero at a rate proportional to  $1/\log n$ , the total MST weight grows approximately linearly in  $n$ . Empirically, we observe

$$f(n) \approx 0.035n.$$

**Geometric graphs (dimensions 2–4).** For  $n$  random points uniformly distributed in  $[0, 1]^d$ , the typical nearest-neighbor distance scales on the order of

$$n^{-1/d}.$$

An MST contains  $n - 1$  edges, and most edges connect nearby points at roughly this scale. Therefore, the total MST weight scales as

$$(n - 1) \cdot n^{-1/d} = \Theta(n^{(d-1)/d}).$$

This matches the empirical behavior:

$$\sqrt{n} \quad (d = 2), \quad n^{2/3} \quad (d = 3), \quad n^{3/4} \quad (d = 4).$$

The experimental constants remain stable as  $n$  increases, indicating convergence toward the asymptotic growth rate.