

# Random Graph Generation

Properties:

- All vertices have an equal chance to be connected to any other vertex
- Any possible graph can be achieved through the algorithm
- Has to be connected

Graph generation consists of two main parts, I will examine each in a separate section. First challenge of random graph generation is the ability to create a random spanning tree which ensures that all of the vertices are connected. Second challenge is adding elements to the spanning tree to the desired density.

## Random Spanning Tree Generation

Inductive building

To ensure that the graph stays connected, we start with a connected graph of size 1 and each subsequent vertex is connected to any part of the graph. We can show by induction that this results in a connected graph.

An easy way of doing this is to start adding elements in a sequence. Any vertices which have been already added can be chosen as the opposite vertex to form an edge.

This method is great to ensure that all vertices are connected. However there are issues with distribution of edges. Since vertices which are added sooner have a higher chance of being connected to, the first 2 properties are violated. Further I will expand on this algorithm.

Chain Building

This is a very efficient way of building a spanning tree which ensures that any vertex has nearly the same probability to be connected to any other vertex. The method is to create an array of all the vertices, shuffle the array and the sequence of the vertices in the array now form a random path through the list which traces out a spanning tree.

This method is however violating the second property, because all graphs where any vertex has a degree  $> 2$  is impossible to achieve.

Inductive building with equal probabilities

This method is very similar to the inductive building, except it favors the last item added to the graph in order to spread out the number of chances each vertex gets to get picked. This is the algorithm I ended up implementing and using to build a spanning tree.

Below is a visualization of my reasoning to calculate the probability using

`random.randint()` linear distribution. Suppose all elements in the array except a last one

had an equal chance to be picked. Since the first element added needs to have a chance to be picked every iteration. The chance each element needs to have is the same as the first element. So when adding  $n$ th element, pick from sequence

$1, 2, \dots, (n-2), (n-1), (n-1), \dots, (n-1)$  where  $(n-1)$  is repeated  $n-1$  times a random element.

```
ADD 2
1 <- pick random element from this sequence

ADD 3
1
122 <- pick random element from this sequence

ADD 4
12
12333 <- pick random element from this sequence

ADD 5
123
123
1234444 <- pick random element from this sequence

ADD 6
1234
1234
1234
123455555 <- pick random element from this sequence
```

This can be implemented in this function:

```
def random_vertex_inductive(n):
    x = random.randint(1, 2*n-1)
    if x < n:
        return x
    return n
```

## Adding Edges

### Random vertex pairs

We begin by having a list of all vertices. After our spanning graph was created we can start adding edges to the graph. This is done by picking 2 random indexes which correspond to 2 random vertices from the vertex list. We try to add them to the graph, which might fail if the

vertices are the same or the graph already has that edge added to it. This initially is not an issue as we can just pick another set of indexes and most likely we would encounter a valid edge to add in expected constant time. This approach starts to be slowed down significantly when we start reaching dense graphs. The probability becomes astronomical for finding the last vertex to add, completing a fully connected graph would be nearly impossible, since the chance to get the last edge is  $n^2/2$ , we would need to try roughly  $n^2/4$  combinations.

This would be unusable, so if we would want to improve on the time it takes to find a correct vertex we could implement linear probing, were we simply advance to the next index.

```
def rollIndexes(indexes, n):
    i1, i2 = indexes
    i1 = (i1+1)%n
    if i1==0:
        i2 = (i2+1)%n
    return i1, i2
```

However this might result in clusters which would mean that the first property is violated.

*Aside: this method could be used by first creating a fully connected graph and then removing edges, while keeping the original spanning tree intact*

## Generating a list of all possible edges

We first generate a list of tuples of 2 vertices, where each tuple represents an edge. We have to make sure not to allow for duplicate edges since order in tuples matters but we are trying to build an undirected graph.

```
def generate_full_connections(n):
    conns = []
    for i in range(n-1):
        for j in range(i+1, n):
            conns.append((i, j))
    random.shuffle(conns)
    return conns
```

I store this list and also keep an index which advances through the list and I store that one as well in the Graph Generator object. This allows to add more edges if needed or remove edges if needed. Of course I need to keep checking if the edge I am about to remove was not in the spanning tree which I keep as a set. Or in case of adding edges I need to check if the edge is not already in the graph from the spanning tree. These are okay checks as they take constant time per edge and they appear overall only  $n - 1$  times at max.

I chose this technique to generate all graphs. Even though the random graph generation is faster when it comes to dense extremes (very dense for random removal or very sparse for random addition) But the random removal is very close in time complexity to this method.

Both methods would generate a fully connected graph which is already  $O(n^2)$  which would be the main complexity regardless of how many edges we add or remove later. And generating sparse graphs might take  $O(n^2)$  for this technique however since it is a sparse graph the creation is quite fast either way.

## Weights

Weight is the third random factor in graph creation. It might seem insignificant at first but further analysis of the prim's algorithm with a binary heap shows that the range of values we allow the weight to take on is going to determine the efficiency of our algorithm to the point where we might get  $O(m + C * \log_2 n)$  where  $C \in \mathbb{R}$ .  $C$  might end up being dependent on the weight range we allow.

---

## Prim's Implementation

I implement the prim's algorithm using pseudocode from the lecture notes. Both adaptable priority queues are being used by identical code. Polymorphism is used here to achieve similar testing conditions

```
def prim(self, APQ_type):
    pq = APQUnsortedList()
    if APQ_type == "heap":
        pq = APQBinaryHeap()
    locs = {}
    for vertex in self.vertices():
        element = pq.add(float("inf"), (vertex, None))
        locs[vertex] = element
    tree = []
    while pq.length() > 0:
        value = pq.remove_min()[1]
        vertex, entry_edge = value
        locs.pop(vertex)
        if entry_edge:
            tree.append(entry_edge)
        for edge in self.get_edges(vertex):
            opposite_vertex = edge.opposite(vertex)
            if locs.get(opposite_vertex):
                cost = edge.weight
                element = locs[opposite_vertex]
                if cost < pq.get_key(element):
                    element.value = (opposite_vertex, edge)
                    pq.update_key(element, cost)
    return tree
```

Quick test to see if the algorithm failed is the following

```
assert len(tree)==(n-1)
```

However this test does not prove that our algorithm is working properly. To check the working of the algorithm I ran it on some smaller samples and then on larger graphs, pasted the results into <https://edotor.net/> to check if the graph was indeed a spanning tree and preferred lower value edges.

For this testing there is a function in `graph_test.py` file which prints out the formatted graph to be checked together with the spanning tree.

## Unsorted List

Implementing the APQ using an unsorted list is quite straightforward. We are aiming for `update_key()` and `remove()` to be  $O(1)$  and `remove_min()` to be  $O(n)$ .

We achieve the constant access and update time by not keeping any kind of structure inside of the underlying array and storing the index in the element objects themselves (These need to be updated anytime they would change position). The user of this class has to manage the element objects themselves, since they need to be passed into methods like `update_key()` and `remove()`.

Linear time for removing the minimum element is done by searching the whole list and removing the element which had the highest priority.

This implementation is efficient for algorithms where updating and removing from the APQ is preferred or more frequent than removing the minimum.

The complexity of this algorithm is  $O(n^2 + m + n)$ , and since for a fully connected graph  $m = \frac{n*(n-1)}{2}$  we can see that the complexity is  $O(n^2)$ .

## Binary Heap

Similarly the binary heap is storing elements which keep their own index inside the array which represents the binary heap. However this time, when we update or remove we are forced to rebalance the heap, which takes  $O(\log n)$ . Same complexity applies to removing a minimum element.

When applied to the prim's algorithm, we should try and avoid performing any actions on the APQ as all of them have the disadvantage of rebalancing the heap. It is also worth noting that for smaller graphs the cost of rebalancing is proportionally higher than for larger ones where the logarithmic growth creates an advantage over for example linear growth.

The time complexity of prim's when using this data structure is  $O(n \log n + n \log n + m \log n)$  and since generally  $m > n$  the complexity becomes  $O(m \log n)$ . This means that the performance of this algorithm is heavily dependent on the density of the graph.

---

## Analysis

I will analyse the runtime of the Prim's Algorithm with basic runtime complexities and check the predictions practically. After collection of data I will analyse the trends and improve the theoretical knowledge based on the data.

## Predictions

From the analysis of runtime complexities we can notice that heap should perform better on sparse graphs while the unsorted list should overtake it the denser the graph is.

More specifically at  $m = \frac{n^2}{\log 2}$ .

We can use 3d graphing tools like [Desmos](#) to help us plot the the correlation between vertices, edges and the runtime.

When interrogating further, we can take the limit of the ratio of edges where Heap would perform better to the total number of vertices.

$$\lim_{n \rightarrow \infty} \left( \frac{\frac{n^2}{\log n} - (n-1)}{\frac{(n-1)n}{2} - (n-1)} \right)$$

This can be simplified down to:

$$\lim_{n \rightarrow \infty} \left( \frac{2n^2}{n^2 \log n} \right) = \lim_{n \rightarrow \infty} \left( \frac{2}{\log n} \right) = 0$$

We can see that the ratio of possible edges on which the unsorted list should be better is growing faster than heap. However we will never reach a point where the unsorted list would perform better on a spanning tree.

Suppose:

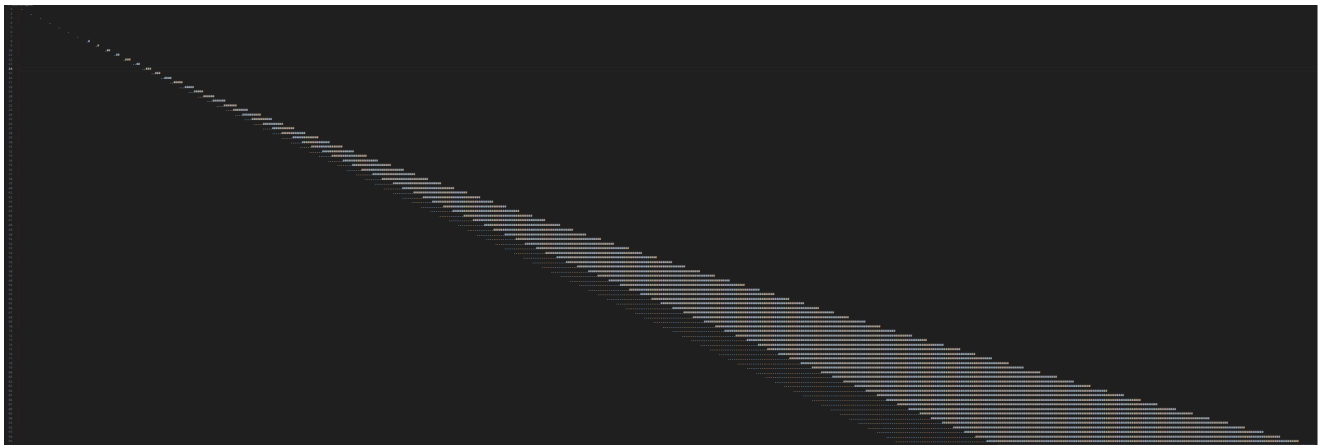
$$\exists N \in \mathbb{N} s.t. \forall n > N \wedge (n) \log n > n^2$$

Then follows:

$$\log n > n$$

Which is a contradiction, thus heap should always be better on sparse graphs, however we know that unsorted list should start taking over larger portions of the list.

*To visualize this, I have created a simple graphing function which plots different graphs as coordinates in a text file and prints a specified character depending on which one is better for that graph*



## Practical Testing

I run simple tests on a range of vertices. For one set of vertices I create a given number of slices from a range of values  $(n - 1; \frac{n*(n-1)}{2})$  endpoints inclusive. These slices give me a rough idea of how the algorithms behave across a spectrum of vertices and densities. Each slice gives me one graph which I measure the performance of Heap Prim and Unsorted List Prim.

### *results*

Vertices=	256	Edges=	255	Heap=0.00000000	Unsorted List=0.00591731
Vertices=	256	Edges=	3494	Heap=0.00904083	Unsorted List=0.00690007
Vertices=	256	Edges=	6733	Heap=0.00675368	Unsorted List=0.00215816
Vertices=	256	Edges=	9972	Heap=0.00769448	Unsorted List=0.01296353
Vertices=	256	Edges=	13211	Heap=0.01459694	Unsorted List=0.01482916
Vertices=	256	Edges=	16450	Heap=0.01349425	Unsorted List=0.02640629
Vertices=	256	Edges=	19688	Heap=0.01801300	Unsorted List=0.01171160
Vertices=	256	Edges=	22926	Heap=0.02000141	Unsorted List=0.01867127
Vertices=	256	Edges=	26164	Heap=0.01984072	Unsorted List=0.01929164
Vertices=	256	Edges=	29402	Heap=0.01780105	Unsorted List=0.02702188
Vertices=	256	Edges=	32640	Heap=0.02942586	Unsorted List=0.02210474
Vertices=	512	Edges=	511	Heap=0.00000000	Unsorted List=0.01873016
Vertices=	512	Edges=	13542	Heap=0.02129960	Unsorted List=0.02161527
Vertices=	512	Edges=	26573	Heap=0.02410030	Unsorted List=0.02698421
Vertices=	512	Edges=	39604	Heap=0.03124380	Unsorted List=0.04017973
Vertices=	512	Edges=	52635	Heap=0.04850054	Unsorted List=0.04719377
Vertices=	512	Edges=	65666	Heap=0.05140662	Unsorted List=0.05749321
Vertices=	512	Edges=	78696	Heap=0.05997229	Unsorted List=0.07684660
Vertices=	512	Edges=	91726	Heap=0.07141519	Unsorted List=0.09285998
Vertices=	512	Edges=	104756	Heap=0.07698655	Unsorted List=0.10003662
Vertices=	512	Edges=	117786	Heap=0.11284113	Unsorted List=0.11713958
Vertices=	512	Edges=	130816	Heap=0.11037731	Unsorted List=0.11667442

Vertices=	1024	Edges=	1023	Heap=0.01284027	Unsorted	List=0.05599833
Vertices=	1024	Edges=	53299	Heap=0.05762053	Unsorted	List=0.09098458
Vertices=	1024	Edges=	105575	Heap=0.09191561	Unsorted	List=0.13834524
Vertices=	1024	Edges=	157851	Heap=0.14905000	Unsorted	List=0.20028639
Vertices=	1024	Edges=	210126	Heap=0.21337247	Unsorted	List=0.26537108
Vertices=	1024	Edges=	262401	Heap=0.26798677	Unsorted	List=0.33139110
Vertices=	1024	Edges=	314676	Heap=0.33599973	Unsorted	List=0.37092638
Vertices=	1024	Edges=	366951	Heap=0.46236229	Unsorted	List=0.47556877
Vertices=	1024	Edges=	419226	Heap=0.51710606	Unsorted	List=0.56461382
Vertices=	1024	Edges=	471501	Heap=0.56991053	Unsorted	List=0.60084701
Vertices=	1024	Edges=	523776	Heap=0.62815285	Unsorted	List=0.68886137
Vertices=	2048	Edges=	2047	Heap=0.02011585	Unsorted	List=0.21645069
Vertices=	2048	Edges=	211456	Heap=0.22594357	Unsorted	List=0.43698263
Vertices=	2048	Edges=	420864	Heap=0.58498549	Unsorted	List=0.70193315
Vertices=	2048	Edges=	630272	Heap=0.74917006	Unsorted	List=0.93990946
Vertices=	2048	Edges=	839680	Heap=1.06798840	Unsorted	List=1.19983196
Vertices=	2048	Edges=	1049088	Heap=1.38133645	Unsorted	List=1.57970238
Vertices=	2048	Edges=	1258496	Heap=1.54128933	Unsorted	List=1.71995401
Vertices=	2048	Edges=	1467904	Heap=1.88867259	Unsorted	List=2.02714539
Vertices=	2048	Edges=	1677312	Heap=2.04961419	Unsorted	List=2.21526217
Vertices=	2048	Edges=	1886720	Heap=2.31984258	Unsorted	List=2.47129464
Vertices=	2048	Edges=	2096128	Heap=2.81836367	Unsorted	List=2.72852230
Vertices=	4096	Edges=	4095	Heap=0.03884435	Unsorted	List=0.89964724
Vertices=	4096	Edges=	842342	Heap=1.07939148	Unsorted	List=1.84855604
Vertices=	4096	Edges=	1680589	Heap=2.28786206	Unsorted	List=3.15204215
Vertices=	4096	Edges=	2518836	Heap=3.19852924	Unsorted	List=3.97246814
Vertices=	4096	Edges=	3357083	Heap=4.39886260	Unsorted	List=5.04938507
Vertices=	4096	Edges=	4195330	Heap=5.38085079	Unsorted	List=6.03921032
Vertices=	4096	Edges=	5033576	Heap=8.17561269	Unsorted	List=8.53456688
Vertices=	4096	Edges=	5871822	Heap=8.00880814	Unsorted	List=8.33761835
Vertices=	4096	Edges=	6710068	Heap=22.72403622	Unsorted	
List=	21.23875761					
Vertices=	4096	Edges=	7548314	Heap=14.27042317	Unsorted	
List=	25.72153616					
Vertices=	4096	Edges=	8386560	Heap=13.26482844	Unsorted	
List=	12.08679676					

## Observations

- We can notice that for sparse graphs, especially the minimum spanning tree, the heap implementation performs much better, this is something we would expect from the complexity analysis.



- We can also see that the jump between vertices with the same density follows  $O(n^2)$  growth. Each step in my testing is two times the previous vertices count and we can see that the runtime jumps by roughly 4 times.
- However the main surprise is that our predictions have not been confirmed. It seems that heap is outperforming the unsorted list even on very dense graphs, sometimes even on fully connected graphs.
  - This should be the direction of our enquiry as deciding which APQ to use is crucial for optimal runtimes.

## Further Analysis

Firstly we need to find an explanation for heap's improved performance when it comes to dense graphs. There has to be some form of optimization in this pseudocode which went unnoticed.

```
prim():
    create an APQ pq, to contain costs and (vertex,edge) pairs
    create an empty dictionary locs for locations of vertices in pq
    for each v in G
        add (∞, (v,None)) into pq and store location in locs[v]
    create an empty list tree, which will be the output
    while pq is not empty
        remove c:(v,e), the minimum element, from pq
        remove v from locs
        if e is not None, append e to tree
        for each edge d incident on v
            w = d's opposite vertex from v
            if w is in locs # and so it is not yet in the tree
                cost = d's cost
                if cost is cheaper than w's entry in pq
                    replace ?:(w,?) in pq with cost: (w,
d)

    return tree
```

We can see that the section of the code which is responsible from the runtime  $O(m \log n)$  is:

```
while pq is not empty
    remove c:(v,e), the minimum element, from pq
    remove v from locs
    if e is not None, append e to tree
    for each edge d incident on v
        w = d's opposite vertex from v
```

```

    if w is in locs # and so it is not yet in the tree
        cost = d's cost
        if cost is cheaper than w's entry in pq
            replace ?:(w,?) in pq with cost: (w, d)

```

Note that the sum of degrees of all vertices is  $2 * m$ , only  $m$  make it through if w is in locs

Here the while loop runs precisely  $n$  times while the inner loop runs for every edge on the vertex  $v$ .

We only update the APQ if cost is cheaper than w's entry in pq. This means that the actual runtime is  $O(m(1 + c \log n))$ . This is interesting because if  $c$  is not a constant but a function of  $m$ , specifically  $\frac{1}{m}$  in some form, the overall runtime might be smaller.

So what is  $c$ ?  $c$  is the probability that the current examined cost is cheaper than the cost in the entry in the APQ. Since it is a probability, it helps to formulate it as a number of possibilities we want: overall number of updates:  $u$  over the overall number of edges  $m$

$$c = \frac{u}{m}$$

Defining  $u$  is tricky since it relies on probability, but I find it helpful to formulate in terms of a single vertex. Each vertex has a degree, this average degree of vertices is  $2 * m/n$ , but only half of the edges are examined, so  $m/n$ . The reason I define the average degree here is because we want to create a set of all possible unique elements the vertex has access to. This is helpful because over the course of the algorithm, as long as it is random, we will keep inspecting elements from this set. We only update the APQ if the element we inspect is lower than the previous one.

If we assume that, on average, the set is split into two by a random item being put inside the APQ, then that creates two equal subsets, where only one will result in the APQ being updated. This splitting of the set continues. So if we assume that a vertex will have  $m/n$  inspections, then the first inspection has a 100% chance to update APQ, second one 50% and so on these probabilities cascade like in a Pascal's triangle where let's say the most left means all were successful updates and the further right you go, the less updates to the APQ have been made.

If we allow for rough estimates as I was not able to find a proper probability: The overall probability for degree  $m/n$  is  $\log n$  roughly.

So now we know that the set can split, now the question is what is size of the set that is being split. Well it is a set of all possible unique elements which can be found on average. First of all there are two limiting factors, if I have 20 weight options, then I can only update a vertex 20 times at max, at best 1 time and on average  $\log_2 20$  times. Similarly I cannot update a vertex more time than it has connections, so  $\log_2 m/n$  or  $\log_2 w$  per vertex updates. Where

$w$  is a weight. We need to pick the smaller one of these because whichever one is smaller will determine what new item cannot be selected anymore.

$$\min\{\log(m/n), \log_2 w\}$$

Putting all of this together;

$$c = \frac{n \cdot \min\{\log(m/n), \log_2 w\}}{m}$$

$$m(1 + c \log n) = m + m \cdot c \log n = m + n \cdot \min\{\log(m/n), \log_2 w\} \cdot \log n$$

$$m + n \cdot \log n \cdot \min\{\log(m/n), \log_2 w\}$$

So if we have the weight set to a small number, like in our test cases where  $w = 20$ , we can view it as a constant and get  $O(m + n \log n)$  as the time complexity for the heap. This would explain why the run times come so close to each other when it comes to Dense Graphs, they are both technically  $O(n^2)$ .

## Further Testing

To ensure my reasoning is correct and the math is giving a good approximation, I measured the total number of times the line: `if cost is cheaper than w's entry in pq`. It was  $m$  and then the total number when the if statement was True has been tested. I have plotted the data with 100 vertices and variable number of edges. We can see that it most likely follows a logarithmic curve.

