

76ljsuxqz

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1 Report

Autors: Mykola Vysotskyi, Kohan Iryna

0. Preparation for testing. In this section we will prepare our data for testing. We just import all necessary libraries and generate graphs for testing.

You can skip this section.

```
[2]: import networkx as nx
      from graph_generator import gnp_random_connected_graph

      from timing import get_timing_data, plot_timing_data
```

Generate graphs for testing.

```
[3]: graphs = tuple(gnp_random_connected_graph(i, 0.1, draw=False) for i in
      ↪range(10, 150, 10))
```

1.1 1. Graph algorithms.

1.1.1 Task 1.1 Kruskal's and Prim's algorithms for minimum spanning tree.

Kruskal's and Prim's algorithms are a greedy algorithms that finds a minimum spanning tree for a connected weighted graph. This means it finds a subset of the edges that forms a tree that includes every vertex, where the total weight of all the edges in the tree is minimized. If the graph is not connected, then it finds a minimum spanning forest (a minimum spanning tree for each connected component).

Kruskal algorithm code.

```
[4]: def kruskal_search(graph: nx.Graph) -> nx.Graph:
      """
      Implements the Kruskal's algorithm to find
      the minimum spanning tree of a graph.
      Returns list[tuple], where tuple is tuple of nodes.

      Parameters
      -----
      graph : nx.Graph
```

Graph to find the minimum spanning tree of.

Returns

nx.Graph : Minimum spanning tree of the graph.

"""

```
graph = graph.to_directed()
edges = sorted(graph.edges(data=True), key = lambda x: x[2]['weight'])

points = [{i} for i in list(graph.nodes())]
result = nx.Graph()
ind1, ind2 = None, None
for edge in edges:
    for point in points:
        if edge[0] in point:
            ind1 = points.index(point)
        if edge[1] in point:
            ind2 = points.index(point)
    if ind1 != ind2:
        points[ind1] = points[ind1].union(points[ind2])
        del points[ind2]

    result.add_edge(edge[0], edge[1], weight=edge[2]['weight'])

return result
```

Prim's algorithm code.

```
[5]: from queue import PriorityQueue

def prim_mst(graph: nx.Graph) -> nx.Graph:
    """
    Implements the Prim's algorithm to find
    the minimum spanning tree of a graph.
    Returns list[tuple], where tuple is tuple of nodes.

    Parameters
    -----
    graph : nx.Graph
        Graph to find the minimum spanning tree of.

    Returns
    -----
    nx.Graph : Minimum spanning tree of the graph.
    """
    graph = graph.to_directed()
    visited = [False] * graph.number_of_nodes()
```

```

visited[0] = True

mst = nx.Graph()
edges = PriorityQueue()

for adj, data in graph.adj[0].items():
    if not visited[adj]:
        edges.put((data["weight"], 0, adj))

while not edges.empty():
    weight, tree_vertex, vertex = edges.get()
    if visited[vertex]:
        continue
    mst.add_edge(tree_vertex, vertex, weight=weight)
    visited[vertex] = True

    for adj, data in graph.adj[vertex].items():
        if not visited[adj]:
            edges.put((data["weight"], vertex, adj))

return mst

```

Let's test our algorithms on our generated graphs.

```

[6]: timing_data = get_timing_data(
    {
        "Prim": (prim_mst, ()),
        "Kruskal": (kruskal_search, ()),
        "Kruskal(networkx)": (nx.tree.maximum_spanning_tree, ()),
    }, graphs)

```

```

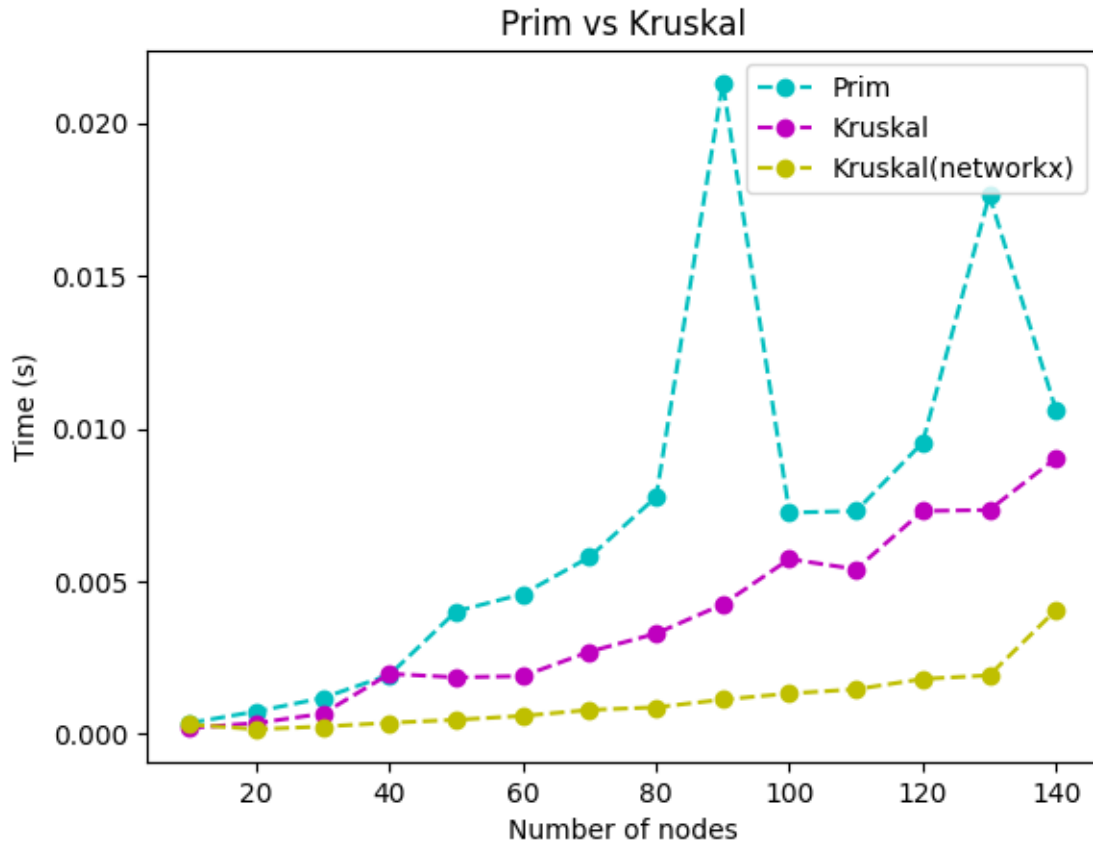
[7]: plot_timing_data(timing_data, title="Prim vs Kruskal", xlabel="Number of ↵
    ↵nodes", ylabel="Time (s)")

```

```

c:\Users\Mykola\Desktop\dm-lab1\timing.py:82: UserWarning: Matplotlib is
currently using module://matplotlib_inline.backend_inline, which is a non-GUI
backend, so cannot show the figure.
    fig.show()

```



As we can see Kruskal's algorithm works faster than Prim's algorithm. In Prim's algorithm implementation we use priority queue, but in Kruskal's algorithm we use sorting. So, we can say that sorting is faster than priority queue.

1.1.2 Task 1.2 Bellman-Ford and Floyd-Warshall algorithms.

Bellman-Ford algorithm code.

```
[4]: def belman_ford_algorithm(graph: nx.Graph, start: int = 0) -> tuple[list[float], bool]:
    """
    Function implements Belman-Ford's algorithm.
    """
    graph = graph.to_directed()
    num_nodes = graph.number_of_nodes()
    edges = graph.edges(data="weight")

    distances = {idx: float("inf") for idx in range(num_nodes)}
    distances[start] = 0
```

```

for _ in range(num_nodes - 1):
    for edge in edges:
        if distances[edge[1]] > (distances[edge[0]] + edge[2]):
            distances[edge[1]] = distances[edge[0]] + edge[2]

for edge in edges:
    if distances[edge[1]] > (distances[edge[0]] + edge[2]):
        return False

return distances, True

```

[13]: *# create wrapper for networkx.single_source_bellman_ford*
to deal with exceptions

```

def bf_nx(graph: nx.Graph) -> tuple[list[float], bool]:
    try:
        res = nx.single_source_bellman_ford(graph, 0)
    except nx.NetworkXUnbounded:
        return False

    return res, True

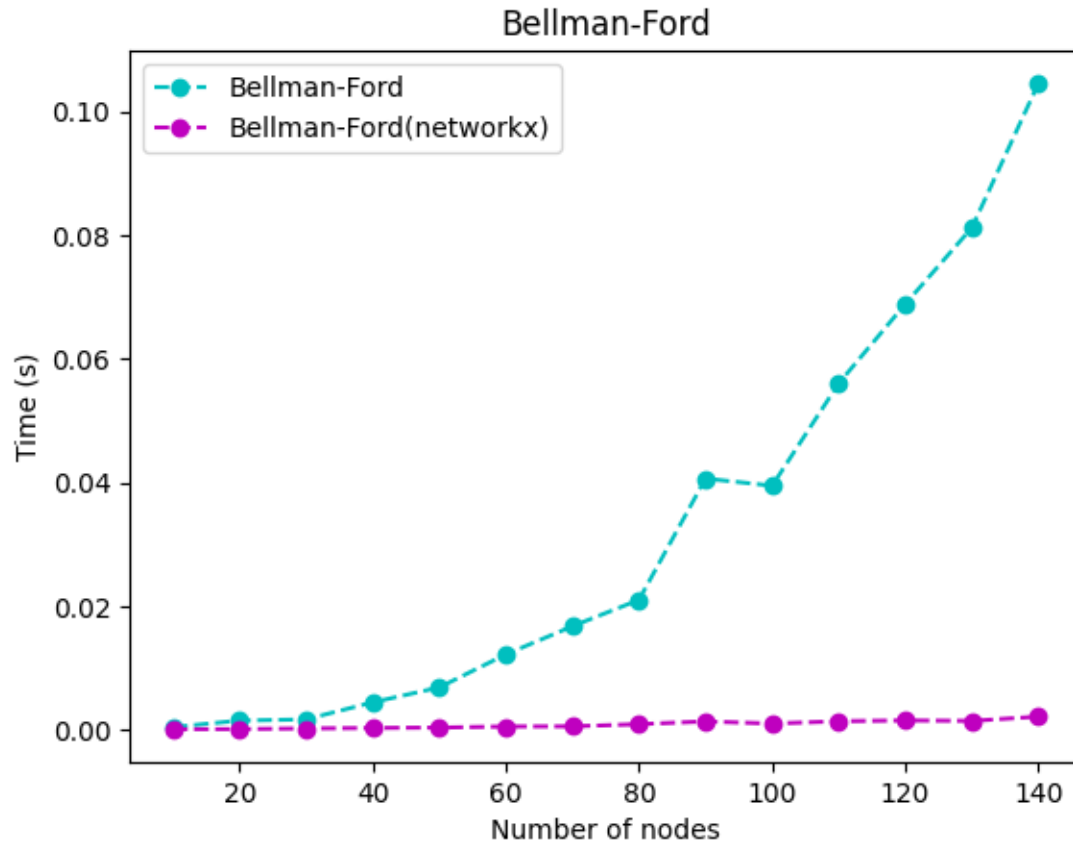
```

```

timing_data = get_timing_data(
{
    "Bellman-Ford": (belman_ford_algorithm, ()),
    "Bellman-Ford(networkx)": (bf_nx, ()),
}, graphs)

```

[14]: `plot_timing_data(timing_data, title="Bellman-Ford", xlabel="Number of nodes",
↪ ylabel="Time (s)")`



Floyd-Warshall algorithm code.

```
[8]: def have_negative_cycle(matrix: list[list[float]]) -> bool:
    """
    Function checks for negative cycle in matrix.
    """
    for idx, _ in enumerate(matrix):
        if matrix[idx][idx] < 0:
            return True

    return False

def floyd_warshall_algorithm(graph: nx.Graph) -> list[list[float]]:
    """
    Function implements Floyd-Warshall's algorithm.
    If there is a negative cycle in graph, function returns None.
    """
    graph = graph.to_directed()
    num_nodes = graph.number_of_nodes()
    edges = graph.edges(data="weight")
```

```

matrix = [[float("inf") for _ in range(num_nodes)] for _ in
↪range(num_nodes)]

for idx in range(num_nodes):
    matrix[idx][idx] = 0

for edge in edges:
    matrix[edge[0]][edge[1]] = edge[2]

for k in range(num_nodes):
    for i in range(num_nodes):
        for j in range(num_nodes):
            if matrix[i][j] > matrix[i][k] + matrix[k][j]:
                matrix[i][j] = matrix[i][k] + matrix[k][j]

if have_negative_cycle(matrix):
    return None

return matrix

```

```

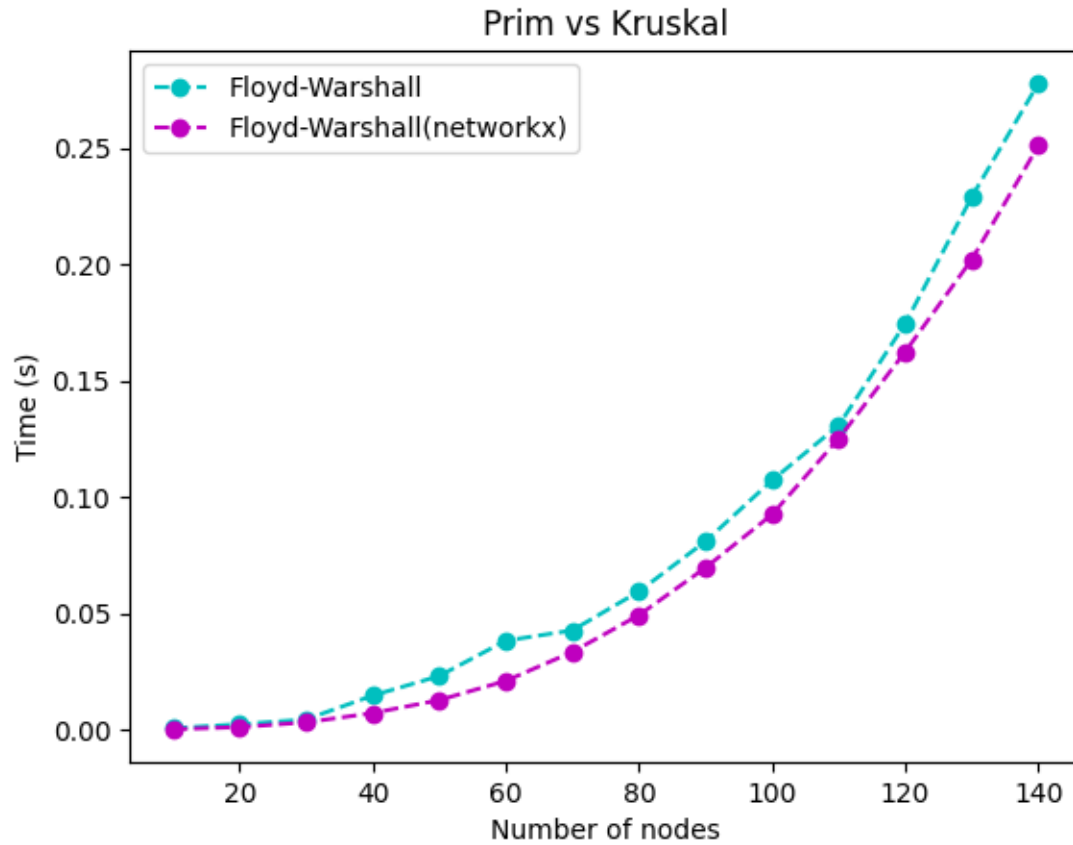
[9]: timing_data = get_timing_data(
    {
        "Floyd-Warshall": (floyd_warshall_algorithm, ()),
        "Floyd-Warshall(networkx)": (nx.floyd_warshall, ())
    }, graphs)

```

```

[10]: plot_timing_data(timing_data, title="Prim vs Kruskal", xlabel="Number of
↪nodes", ylabel="Time (s)")

```



As we can see both algorithms work the same time.

1.2 Task 2. Decision tree classifier.

This task is completed in the file [DecisionTreeClassifier](#). But also we implemented another decision tree algorithm called **ID3**. You can find it in the file [ID3](#).