Simulated annealing minimum spanning tree

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import random
import math
from collections import defaultdict
class Graph:
  def __init__(self):
    self.edges = defaultdict(list)
  def add_edge(self, u, v, weight):
    self.edges[u].append((v, weight))
    self.edges[v].append((u, weight)) # Undirected graph
  def get_edges(self):
    return [(u, v, weight) for u in self.edges for v, weight in self.edges[u] if u < v]
def random_spanning_tree(graph):
  nodes = list(graph.edges.keys())
  random.shuffle(nodes)
  tree_edges = set()
  selected = {nodes[0]}
  while len(selected) < len(nodes):
    u = random.choice(list(selected))
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candidates = [(v, weight) for v, weight in graph.edges[u] if v not in selected]
    if candidates:
       v, weight = random.choice(candidates)
       tree_edges.add((u, v, weight))
       selected.add(v)
  return tree_edges
def energy(tree):
  return sum(weight for u, v, weight in tree)
def generate_neighbor(tree, graph):
  tree_list = list(tree)
  if len(tree_list) < 2:</pre>
    return tree
  # Select a random edge to remove
  u, v, weight = random.choice(tree_list)
  new_tree = tree - {(u, v, weight)}
  # Find a new edge to add
  candidates = [(x, w) \text{ for } x, w \text{ in graph.edges}[u] \text{ if } (x, u, w) \text{ not in tree and } (u, x, w) \text{ not in tree}]
  if not candidates:
    # If no candidates are available, return the original tree
```

return tree

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new_v, new_weight = random.choice(candidates)
  # Add the new edge and check for cycles
  new_tree.add((u, new_v, new_weight))
  # Ensure the new tree is valid (could add a check here if necessary)
  return new_tree
def simulated_annealing(graph):
 T = 1.0 # Initial temperature
 final_temperature = 0.001
  cooling_factor = 0.95
  current_solution = random_spanning_tree(graph)
  best_solution = current_solution
  while T > final_temperature:
    for _ in range(100): # Number of iterations at current temperature
      neighbor = generate_neighbor(current_solution, graph)
      current_energy = energy(current_solution)
      neighbor_energy = energy(neighbor)
      if neighbor_energy < current_energy:</pre>
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current_solution = neighbor
       else:
         acceptance_probability = math.exp((current_energy - neighbor_energy) / T)
         if random.random() < acceptance_probability:</pre>
           current_solution = neighbor
      if energy(current_solution) < energy(best_solution):</pre>
         best_solution = current_solution
    T *= cooling_factor
  return best_solution
# Example usage:
if __name__ == "__main__":
  graph = Graph()
  edges = [(0, 1, 4), (0, 2, 1), (1, 2, 2), (1, 3, 5), (2, 3, 3)]
  for u, v, weight in edges:
    graph.add_edge(u, v, weight)
  mst = simulated_annealing(graph)
  print("Edges in the Minimum Spanning Tree:")
  for u, v, weight in mst:
    print(f"{u} -- {v} (weight: {weight})")
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print("Total weight:", energy(mst))

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Edges in the Minimum Spanning Tree:
2 -- 0 (weight: 1)
2 -- 3 (weight: 3)
2 -- 1 (weight: 2)
Total weight: 6
```