```
import random
def generate chromosome(length):
   return [random.randint(0, 1) for _ in range(length)]
def generate_population(size, chromosome_length):
   return [generate_chromosome(chromosome_length) for _ in range(size)]
def fitness(chromosome):
   # Implement your fitness function here
   # This is a placeholder example, replace it with your actual fitness calculation
   return sum(chromosome)
def selection(population, fitnesses):
   # Implement your selection mechanism here
   # This is a simple roulette wheel selection example
   total_fitness = sum(fitnesses)
   probabilities = [f / total fitness for f in fitnesses]
   return random.choices(population, weights=probabilities, k=2)
def crossover(parent1, parent2):
   crossover_point = random.randint(1, len(parent1) - 1)
   offspring1 = parent1[:crossover_point] + parent2[crossover_point:]
   offspring2 = parent2[:crossover_point] + parent1[crossover_point:]
   return offspring1, offspring2
def mutation(chromosome, mutation_rate):
   for i in range(len(chromosome)):
       if random.random() < mutation rate:</pre>
           chromosome[i] = 1 - chromosome[i]
   return chromosome
def genetic_algorithm(population_size, chromosome_length, generations, mutation_rate):
   population = generate_population(population_size, chromosome_length)
   for generation in range(generations):
       fitnesses = [fitness(chromosome) for chromosome in population]
       new_population = []
       for _ in range(population_size // 2):
           parent1, parent2 = selection(population, fitnesses)
          offspring1, offspring2 = crossover(parent1, parent2)
           offspring1 = mutation(offspring1, mutation_rate)
          offspring2 = mutation(offspring2, mutation_rate)
          new_population.extend([offspring1, offspring2])
       population = new_population
       best_fitness = max(fitnesses)
       best_chromosome = population[fitnesses.index(best_fitness)]
       print(f"Generation {generation+1}: Best fitness = {best_fitness}, Chromosome = {best_chromosome}")
   return best_chromosome
# Example usage:
population_size = 100
chromosome length = 20
generations = 100
mutation_rate = 0.01
\verb|best_solution| = \verb|genetic_algorithm| (population_size, chromosome_length, generations, mutation_rate)| \\
print("Final best solution:", best_solution)
Fraction 1: Best fitness = 15, Chromosome = [0, 1, 1, 1, 0, 1, 0, 1, 1, 1, 0, 0, 1, 1, 1, 0, 1, 1, 1, 1]
    Generation 2: Best fitness = 16, Chromosome = [1, 0, 1, 0, 1, 0, 1, 0, 1, 1, 0, 1, 0, 1, 1, 0, 1, 0, 0]
    Generation 3: Best fitness = 16, Chromosome = [0, 0, 0, 1, 1, 1, 0, 0, 0, 1, 0, 1, 0, 1, 0, 1, 0, 0, 0, 0, 1]
    Generation 4: Best fitness = 16, Chromosome = [1, 1, 0, 0, 1, 1, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 1, 1, 1, 0]
    Generation 5: Best fitness = 16, Chromosome = [1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 0, 1, 0, 0, 1, 1, 1, 0, 0]
    Generation 6: Best fitness = 16, Chromosome = [0, 1, 1, 1, 0, 1, 0, 0, 1, 1, 0, 0, 1, 1, 0, 1, 1, 1]
    Generation 7: Best fitness = 16, Chromosome = [0, 1, 1, 0, 1, 0, 0, 1, 1, 1, 0, 0, 1, 1, 1, 0, 1, 1, 1, 0]
    Generation 9: Best fitness = 17, Chromosome = [1, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1]
    Generation 10: Best fitness = 18, Chromosome = [0, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1, 1, 1, 1, 0, 0, 0, 1, 1, 1, 1]
    Generation 11: Best fitness = 19, Chromosome = [0, 1, 1, 1, 0, 1, 1, 1, 0, 0, 1, 1, 1, 1, 1, 1, 1, 1, 0, 1, 1, 0]
     \text{Generation 13: Best fitness = 18, Chromosome = [0, 1, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1] } 
     \text{Generation 14: Best fitness = 19, Chromosome = [0, 1, 1, 1, 0, 1, 1, 1, 0, 0, 1, 0, 1, 1, 1, 1, 1, 1, 0, 1, 1] } 
    Generation 15: Best fitness = 19, Chromosome = [1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 1, 1, 1]
    Generation 16: Best fitness = 19, Chromosome = [1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1]
    Generation 19: Best fitness = 19, Chromosome = [1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 1, 0, 1, 1, 0, 1, 1]
     \text{Generation 20: Best fitness = 18, Chromosome = [1, 1, 1, 0, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1, 0, 0, 1, 1] }
```

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Generation 22: Best fitness = 18, Chromosome = [1, 1, 0, 0, 1, 1, 0, 1, 0, 1, 1, 1, 1, 1, 0, 1, 1, 0, 0, 1, 0]
   Generation 24: Best fitness = 19, Chromosome = [0, 1, 1, 0, 1, 0, 0, 1, 1, 1, 1, 0, 0, 1, 1, 1, 1, 1, 1, 1]
   Generation 25: Best fitness = 19, Chromosome = [0, 1, 1, 1, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 0, 1, 0, 1, 1]
   Generation 26: Best fitness = 18, Chromosome = [0, 1, 1, 1, 1, 0, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 0, 1, 1]
   Generation 31: Best fitness = 19, Chromosome = [0, 1, 1, 1, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1]
   Generation 32: Best fitness = 19, Chromosome = [1, 1, 0, 0, 1, 1, 1, 1, 1, 0, 1, 0, 1, 1, 1, 0, 1, 0, 1, 0]
   Generation 34: Best fitness = 19, Chromosome = [1, 0, 1, 1, 0, 1, 1, 0, 0, 1, 0, 0, 1, 0, 1, 1, 1, 1, 1, 1]
   Generation 35: Best fitness = 19, Chromosome = [1, 1, 1, 1, 0, 1, 1, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 0, 0, 0, 1]
   Generation 36: Best fitness = 20, Chromosome = [1, 1, 1, 1, 0, 1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 0, 1]
   Generation 37: Best fitness = 19, Chromosome = [1, 1, 1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 1, 1, 1]
   Generation 39: Best fitness = 19, Chromosome = [0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 1, 0, 1, 1, 1, 1, 1, 1]
   Generation 45: Best fitness = 20, Chromosome = [1, 1, 1, 1, 1, 0, 0, 0, 1, 0, 1, 0, 1, 1, 1, 1, 1, 0, 0, 1]
   Generation 47: Best fitness = 20, Chromosome = [1, 1, 1, 1, 1, 1, 0, 1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1]
   Generation 50: Best fitness = 20, Chromosome = [1, 0, 1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 0, 1, 1, 1, 0, 1, 1]
   Generation 52: Best fitness = 20, Chromosome = [1, 0, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 0, 1, 1, 1, 1, 0, 0, 1]
   Generation 55: Best fitness = 20, Chromosome = [1, 1, 1, 0, 1, 1, 1, 1, 1, 0, 1, 1, 0, 1, 0, 1, 1, 0, 1, 1]
   Generation 57: Best fitness = 20, Chromosome = [1, 0, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1, 1, 0, 1, 0]
   Generation 58: Best fitness = 20, Chromosome = [1, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 1, 1, 0, 1, 1, 1]
import random
import numpy as np
class Particle:
  def init (self, dimensions):
     self.position = [random.uniform(-10, 10) for _ in range(dimensions)]
     self.velocity = [random.uniform(-1, 1) for _ in range(dimensions)]
     self.pbest_position = self.position.copy()
     self.pbest_value = float('inf')
class PSO:
  def __init__(self, dimensions, swarm_size, iterations, c1, c2, w):
     self.dimensions = dimensions
     self.swarm_size = swarm_size
     self.iterations = iterations
     self.c1 = c1 # Cognitive coefficient
     self.c2 = c2 # Social coefficient
     self.w = w # Inertia weight
     self.swarm = [Particle(dimensions) for _ in range(swarm_size)]
     # Initialize gbest based on initial population fitness
     self.gbest_value = float('inf')
     for particle in self.swarm:
       fitness = self.evaluate(particle.position)
       if fitness < self.gbest_value:</pre>
          self.gbest_value = fitness
          self.gbest_position = particle.position.copy()
  def evaluate(self, position):
     # Replace this with your specific objective function
     return sum(x**2 \text{ for } x \text{ in position})
  def update_velocity(self, particle):
     for i in range(self.dimensions):
       r1 = random.random()
       r2 = random.random()
       particle.velocity[i] = (self.w * particle.velocity[i]) + \
                     (self.c1 * r1 * (particle.pbest_position[i] - particle.position[i])) + \
(self.c2 * r2 * (self.gbest_position[i] - particle.position[i]))
  def update_position(self, particle):
     for i in range(self.dimensions):
       particle.position[i] += particle.velocity[i]
  def update_pbest(self, particle):
```

```
fitness = self.evaluate(particle.position)
        if fitness < particle.pbest value:</pre>
            particle.pbest_value = fitness
            particle.pbest_position = particle.position.copy()
    def update_gbest(self):
        for particle in self.swarm:
            fitness = self.evaluate(particle.position)
            if fitness < self.gbest_value:</pre>
                self.gbest value = fitness
                self.gbest_position = particle.position.copy()
    def optimize(self):
        for _ in range(self.iterations):
            for particle in self.swarm:
                self.update_velocity(particle)
                self.update_position(particle)
                self.update_pbest(particle)
            self.update_gbest()
        return self.gbest_position, self.gbest_value
# Example usage:
dimensions = 10
swarm_size = 50
iterations = 100
c1 = 2.0
c2 = 2.0
W = 0.7
pso = PSO(dimensions, swarm_size, iterations, c1, c2, w)
best_position, best_value = pso.optimize()
print("Best position:", best_position)
print("Best value:", best_value)
    Best position: [1.5804132675985314, 0.6945689028300615, -0.2134268303825733, -0.8610857084066699, -0.30811488979935336, 0.359736818
     Best value: 5.275664017890163
import numpy as np
import random
class AntHillOptimization:
    def __init__(self, distance_matrix, n_ants, n_iterations, alpha=1, beta=5, rho=0.5, q=100):
        self.distance matrix = distance matrix # Distance matrix
        self.num_cities = len(distance_matrix) # Number of cities
        self.n ants = n ants # Number of ants
        self.n\_iterations = n\_iterations # Number of iterations
        self.alpha = alpha # Influence of pheromone
        self.beta = beta # Influence of distance
        self.rho = rho # Evaporation rate of pheromone
        self.q = q # Constant for pheromone update
        self.pheromone = np.ones((self.num_cities, self.num_cities)) # Initial pheromone levels
    def _get_probabilities(self, current_city, visited_cities):
        # Calculate the probabilities for each possible next city based on pheromone and distance
        probabilities = []
        total = 0.0
        for i in range(self.num cities):
            if i not in visited cities:
                pheromone = self.pheromone[current_city][i] ** self.alpha
                distance = float(self.distance_matrix[current_city][i]) ** -self.beta # Cast to float to allow negative powers
                probabilities.append(pheromone * distance)
                total += pheromone * distance
            else:
                probabilities.append(0)
        # Normalize probabilities
        probabilities = [prob / total for prob in probabilities]
        return probabilities
    def construct solution(self):
        visited_cities = [random.randint(0, self.num_cities - 1)] # Start from a random city
        while len(visited_cities) < self.num_cities:</pre>
            current_city = visited_cities[-1]
            probabilities = self._get_probabilities(current_city, visited_cities)
            next_city = np.random.choice(self.num_cities, p=probabilities)
            visited_cities.append(next_city)
        visited_cities.append(visited_cities[0]) # Return to the starting city
        return visited cities
```

```
def calculate total distance(self, tour):
        # Calculate the total distance for a given tour
        total_distance = 0
        for i in range(len(tour) - 1):
           total_distance += self.distance_matrix[tour[i]][tour[i + 1]]
        return total_distance
    def _update_pheromone(self, all_solutions, all_distances):
        # Evaporate pheromone levels
        self.pheromone *= (1 - self.rho)
        \# Deposit pheromone based on the quality of solutions
        for i in range(self.n_ants):
           tour = all solutions[i]
            tour_distance = all_distances[i]
            pheromone_deposit = self.q / tour_distance # Deposit pheromone inversely proportional to the distance
            for i in range(len(tour) - 1):
                self.pheromone[tour[i]][tour[i + 1]] += pheromone_deposit
                self.pheromone[tour[i + 1]][tour[i]] += pheromone_deposit # Since the graph is undirected
    def run(self):
        best tour = None
        best_distance = float('inf')
        for iteration in range(self.n_iterations):
            all_solutions = []
            all_distances = []
            for _ in range(self.n_ants):
                solution = self._construct_solution()
                total_distance = self._calculate_total_distance(solution)
                all solutions.append(solution)
                all_distances.append(total_distance)
                if total_distance < best_distance:</pre>
                    best_tour = solution
                    best_distance = total_distance
            # Update pheromone based on the ants' solutions
            self. update pheromone(all solutions, all distances)
            print(f"Iteration {iteration + 1}/{self.n_iterations}, Best Distance: {best_distance}")
        return best_tour, best_distance
# Example usage
if __name__ == "__main__":
    # Example distance matrix for 5 cities
    distance_matrix = np.array([
       [0, 10, 15, 20, 25],
        [10, 0, 35, 25, 30],
       [15, 35, 0, 30, 5],
        [20, 25, 30, 0, 15],
        [25, 30, 5, 15, 0]
    1)
    # Parameters: n_ants, n_iterations, alpha, beta, rho, q
    aco = AntHillOptimization(distance_matrix, n_ants=10, n_iterations=100, alpha=1, beta=5, rho=0.5, q=100)
    best_tour, best_distance = aco.run()
    print(f"Best tour: {best_tour}")
    print(f"Best distance: {best_distance}")

→ Iteration 1/100, Best Distance: 70
     Iteration 2/100, Best Distance: 70
     Iteration 3/100, Best Distance: 70
     Iteration 4/100, Best Distance: 70
     Iteration 5/100, Best Distance: 70
     Iteration 6/100, Best Distance: 70
     Iteration 7/100, Best Distance: 70
     Iteration 8/100, Best Distance: 70
     Iteration 9/100, Best Distance: 70
     Iteration 10/100, Best Distance: 70
     Iteration 11/100, Best Distance: 70
    Iteration 12/100, Best Distance: 70
Iteration 13/100, Best Distance: 70
     Iteration 14/100, Best Distance: 70
     Iteration 15/100, Best Distance: 70
     Iteration 16/100, Best Distance: 70
     Iteration 17/100, Best Distance: 70
```

```
Iteration 18/100, Best Distance: 70
     Iteration 19/100, Best Distance: 70
     Iteration 20/100, Best Distance: 70
     Iteration 21/100, Best Distance: 70
     Iteration 22/100, Best Distance: 70
     Iteration 23/100, Best Distance: 70
     Iteration 24/100, Best Distance: 70
     Iteration 25/100, Best Distance: 70
     Iteration 26/100, Best Distance: 70
     Iteration 27/100, Best Distance: 70
     Iteration 28/100, Best Distance: 70
     Iteration 29/100, Best Distance: 70
     Iteration 30/100, Best Distance: 70
     Iteration 31/100, Best Distance: 70
     Iteration 32/100, Best Distance: 70
     Iteration 33/100, Best Distance: 70
     Iteration 34/100, Best Distance: 70
     Iteration 35/100, Best Distance: 70
     Iteration 36/100, Best Distance: 70
     Iteration 37/100, Best Distance: 70
     Iteration 38/100, Best Distance: 70
     Iteration 39/100, Best Distance: 70
     Iteration 40/100, Best Distance: 70
     Iteration 41/100, Best Distance: 70
     Iteration 42/100, Best Distance: 70
     Iteration 43/100, Best Distance: 70
     Iteration 44/100, Best Distance: 70
     Iteration 45/100, Best Distance: 70
     Iteration 46/100, Best Distance: 70
     Iteration 47/100, Best Distance: 70
     Iteration 48/100, Best Distance: 70
     Iteration 49/100, Best Distance: 70
     Iteration 50/100, Best Distance: 70
     Iteration 51/100, Best Distance: 70
     Iteration 52/100, Best Distance: 70
     Iteration 53/100, Best Distance: 70
     Iteration 54/100, Best Distance: 70
     Iteration 55/100, Best Distance: 70
     Iteration 56/100, Best Distance: 70
     Iteration 57/100, Best Distance: 70
     Iteration 58/100, Best Distance: 70
import numpy as np
import random
# Cuckoo Search for TSP
class CuckooSearch:
    def __init__(self, distance_matrix, n_nests, n_iterations, alpha=0.01, beta=1.5, pa=0.25):
        self.distance_matrix = distance_matrix # Distance matrix
        self.num_cities = len(distance_matrix) # Number of cities
        self.n_nests = n_nests # Number of nests
        self.n_iterations = n_iterations # Number of iterations
        self.alpha = alpha # Step size (scaling factor for Levy flights)
        self.beta = beta # Levy flight exponent
        self.pa = pa # Discovery rate (probability of abandoning a nest)
        {\tt self.nests = np.array([self.\_generate\_solution() \ for \_in \ range(n\_nests)]) \ \# \ Initial \ solutions \ (nests)}
        self.best_solution = None # Best solution
        self.best_distance = float('inf') # Best distance (fitness)
    def _generate_solution(self):
        # Generate a random solution (random permutation of cities)
        return np.random.permutation(self.num_cities)
    def _calculate_total_distance(self, tour):
        # Calculate the total distance of the tour
        total distance = 0
        for i in range(len(tour) - 1):
            total_distance += self.distance_matrix[tour[i]][tour[i + 1]]
        total_distance += self.distance_matrix[tour[-1]][tour[0]] # Return to the starting city
        return total_distance
    def _levy_flight(self, solution):
        # Perform a Levy flight to create a new candidate solution
        step = np.random.normal(0, 1, size=solution.shape) * np.abs(np.random.normal(0, 1)) ** self.beta
        new_solution = solution + self.alpha * step
        # Ensure the new solution is a valid permutation of cities
        new_solution = np.clip(new_solution, 0, self.num_cities - 1).astype(int)
        # Fix invalid solutions (duplicates, out-of-bound values)
        unique_solution = np.unique(new_solution)
        if len(unique_solution) < self.num_cities:</pre>
            # If the solution contains duplicates, replace them with missing cities
            missing_cities = set(range(self.num_cities)) - set(unique_solution)
            new_solution = np.array(list(unique_solution) + list(missing_cities))
```

```
# Ensure the solution has the correct length and is a valid permutation
       return np.random.permutation(new solution) # Ensure it's a permutation
    def _get_best_nests(self, fitness):
       # Get the best solutions (nests) based on fitness (shortest distance)
       sorted_indices = np.argsort(fitness)
       best_nests = self.nests[sorted_indices[:self.n_nests // 2]]
       return best_nests
    def _abandon_nest(self):
       # Abandon the worst nests with probability pa and generate a new solution
       new_nests = []
       for nest in self.nests:
           if random.random() > self.pa:
              new_nests.append(nest)
            else:
               new_nests.append(self._generate_solution()) # Generate a random new solution
       return np.array(new nests)
    def run(self):
       # Main Cuckoo Search loop
        for iteration in range(self.n_iterations):
           fitness = np.array([self._calculate_total_distance(nest) for nest in self.nests])
           # Update the best solution found so far
           best nest idx = np.argmin(fitness)
           best_solution = self.nests[best_nest_idx]
           best_distance = fitness[best_nest_idx]
           if best_distance < self.best_distance:</pre>
               self.best_solution = best_solution
               self.best_distance = best_distance
           # Perform Levy flights to update nests
           new_nests = []
           for i, nest in enumerate(self.nests):
               new_solution = self._levy_flight(nest)
               new_nests.append(new_solution)
           # Ensure all solutions are valid and consistent in shape
           new nests = np.array(new nests)
           # Replace the worst nests with better ones from Levy flights
           fitness_new = np.array([self._calculate_total_distance(nest) for nest in new_nests])
           best_nests = self._get_best_nests(fitness_new)
           self.nests = np.vstack((best_nests, self._abandon_nest()))
           print(f"Iteration {iteration + 1}/{self.n_iterations}, Best Distance: {self.best_distance}")
        return self.best_solution, self.best_distance
# Example usage
if __name__ == "__main__":
   # Example distance matrix for 5 cities
    distance_matrix = np.array([
       [0, 10, 15, 20, 25],
       [10, 0, 35, 25, 30],
       [15, 35, 0, 30, 5],
       [20, 25, 30, 0, 15],
       [25, 30, 5, 15, 0]
    1)
    # Parameters: n_nests, n_iterations, alpha, beta, pa
    cuckoo = CuckooSearch(distance_matrix, n_nests=10, n_iterations=100, alpha=0.01, beta=1.5, pa=0.25)
    best solution, best distance = cuckoo.run()
    print(f"Best solution: {best_solution}")
   print(f"Best distance: {best_distance}")
Iteration 2/100, Best Distance: 95
     Iteration 3/100, Best Distance: 85
     Iteration 4/100, Best Distance: 85
     Iteration 5/100, Best Distance: 85
     Iteration 6/100, Best Distance: 85
     Iteration 7/100, Best Distance: 85
     Iteration 8/100, Best Distance: 85
     Iteration 9/100, Best Distance: 70
     Iteration 10/100, Best Distance: 70
     Iteration 11/100, Best Distance: 70
```

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Iteration 12/100, Best Distance: 70
     Iteration 13/100, Best Distance: 70
     Iteration 14/100, Best Distance: 70
     Iteration 15/100, Best Distance: 70
     Iteration 16/100, Best Distance: 70
     Iteration 17/100, Best Distance: 70
     Iteration 18/100, Best Distance: 70
     Iteration 19/100, Best Distance: 70
     Iteration 20/100, Best Distance: 70
     Iteration 21/100, Best Distance: 70
     Iteration 22/100, Best Distance: 70
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     Iteration 53/100, Best Distance: 70
     Iteration 54/100, Best Distance: 70
     Iteration 55/100, Best Distance: 70
     Iteration 56/100, Best Distance: 70
     Iteration 57/100, Best Distance: 70
     Thomation FO/100 Boot Dictance: 70
import numpy as no
import random
class GreyWolfOptimization:
    def __init__(self, distance_matrix, n_wolves, n_iterations, alpha=0.5, beta=1.5, delta=1.5):
        self.distance_matrix = distance_matrix # Distance matrix
        self.num_cities = len(distance_matrix) # Number of cities
        self.n_wolves = n_wolves # Number of wolves
        self.n_iterations = n_iterations # Number of iterations
        self.alpha = alpha # Coefficient for alpha wolf
        self.beta = beta # Coefficient for beta wolf
        self.delta = delta # Coefficient for delta wolf
        self.wolves = np.array([self._generate_solution() for _ in range(n_wolves)]) # Initial solutions (wolves)
        self.best_solution = None # Best solution
        self.best_distance = float('inf') # Best distance (fitness)
    def _generate_solution(self):
        # Generate a random solution (random permutation of cities)
        return np.random.permutation(self.num_cities)
    def _calculate_total_distance(self, tour):
        # Calculate the total distance of the tour
        total_distance = 0
        for i in range(len(tour) - 1):
           total_distance += self.distance_matrix[tour[i]][tour[i + 1]]
        total_distance += self.distance_matrix[tour[-1]][tour[0]] # Return to the starting city
        return total_distance
    def _update_wolf_position(self, wolf, A, C, alpha_wolf, beta_wolf, delta_wolf):
        # Update position of the wolf using the formula for alpha, beta, and delta wolves
        new_wolf_position = wolf + A * (alpha_wolf - wolf) + C * (beta_wolf - wolf) + C * (delta_wolf - wolf)
        # Make sure the solution is within the valid range [0, num_cities-1] and is a valid permutation
        new_wolf_position = np.clip(new_wolf_position, 0, self.num_cities - 1).astype(int)
        unique_wolf = np.unique(new_wolf_position)
        if len(unique_wolf) < self.num_cities:</pre>
            missing_cities = set(range(self.num_cities)) - set(unique_wolf)
            new_wolf_position = np.array(list(unique_wolf) + list(missing_cities))
        return np.random.permutation(new_wolf_position) # Ensure it's a valid permutation
```

```
def run(self):
        # Main Grey Wolf Optimization loop
        for iteration in range(self.n_iterations):
            fitness = np.array([self._calculate_total_distance(wolf) for wolf in self.wolves])
            # Update the best solution found so far
            best wolf idx = np.argmin(fitness)
            best_solution = self.wolves[best_wolf_idx]
            best_distance = fitness[best_wolf_idx]
            if best_distance < self.best_distance:</pre>
                self.best_solution = best_solution
                self.best_distance = best_distance
            # Get alpha, beta, and delta wolves
            sorted_indices = np.argsort(fitness)
            alpha wolf = self.wolves[sorted indices[0]]
            beta_wolf = self.wolves[sorted_indices[1]]
            delta_wolf = self.wolves[sorted_indices[2]]
            \mbox{\tt\#} Update positions of the wolves using the GWO formula
            A = 2 * np.random.random(self.wolves.shape) - 1 # Random vector for A
            C = 2 * np.random.random(self.wolves.shape) # Random vector for C
            new wolves = []
            for i, wolf in enumerate(self.wolves):
               new_position = self._update_wolf_position(wolf, A[i], C[i], alpha_wolf, beta_wolf, delta_wolf)
                new_wolves.append(new_position)
            self.wolves = np.array(new_wolves)
            print(f"Iteration {iteration + 1}/{self.n_iterations}, Best Distance: {self.best_distance}")
        return self.best_solution, self.best_distance
# Example usage
if __name__ == "__main__":
   # Example distance matrix for 5 cities
    distance_matrix = np.array([
       [0, 10, 15, 20, 25],
        [10, 0, 35, 25, 30],
        [15, 35, 0, 30, 5],
       [20, 25, 30, 0, 15],
       [25, 30, 5, 15, 0]
    ])
    # Parameters: n_wolves, n_iterations, alpha, beta, delta
    gwo = GreyWolfOptimization(distance_matrix, n_wolves=10, n_iterations=100, alpha=0.5, beta=1.5, delta=1.5)
    best_solution, best_distance = gwo.run()
    print(f"Best solution: {best_solution}")
   print(f"Best distance: {best_distance}")
Iteration 2/100, Best Distance: 70
     Iteration 3/100, Best Distance: 70
     Iteration 4/100, Best Distance: 70
Iteration 5/100, Best Distance: 70
     Iteration 6/100, Best Distance: 70
     Iteration 7/100, Best Distance: 70
     Iteration 8/100, Best Distance: 70
     Iteration 9/100, Best Distance: 70
     Iteration 10/100, Best Distance: 70
     Iteration 11/100, Best Distance: 70
     Iteration 12/100, Best Distance: 70
     Iteration 13/100, Best Distance: 70
     Iteration 14/100, Best Distance: 70
     Iteration 15/100, Best Distance: 70
     Iteration 16/100, Best Distance: 70
     Iteration 17/100, Best Distance: 70
     Iteration 18/100, Best Distance: 70
     Iteration 19/100, Best Distance: 70
     Iteration 20/100, Best Distance: 70
     Iteration 21/100, Best Distance: 70
     Iteration 22/100, Best Distance: 70
     Iteration 23/100, Best Distance: 70
     Iteration 24/100, Best Distance: 70
     Iteration 25/100, Best Distance: 70
     Iteration 26/100, Best Distance: 70
     Iteration 27/100, Best Distance: 70
     Iteration 28/100, Best Distance: 70
     Iteration 29/100, Best Distance: 70
```

```
Iteration 30/100, Best Distance: 70
     Iteration 31/100, Best Distance: 70
     Iteration 32/100, Best Distance: 70
     Iteration 33/100, Best Distance: 70
     Iteration 34/100, Best Distance: 70
     Iteration 35/100, Best Distance: 70
     Iteration 36/100, Best Distance: 70
     Iteration 37/100, Best Distance: 70
     Iteration 38/100, Best Distance: 70
     Iteration 39/100, Best Distance: 70
     Iteration 40/100, Best Distance: 70
     Iteration 41/100, Best Distance: 70
     Iteration 42/100, Best Distance: 70
     Iteration 43/100, Best Distance: 70
     Iteration 44/100, Best Distance: 70
     Iteration 45/100, Best Distance: 70
     Iteration 46/100, Best Distance: 70
     Iteration 47/100, Best Distance: 70
     Iteration 48/100, Best Distance: 70
     Iteration 49/100, Best Distance: 70
     Iteration 50/100, Best Distance: 70
     Iteration 51/100, Best Distance: 70
     Iteration 52/100, Best Distance: 70
     Iteration 53/100, Best Distance: 70
     Iteration 54/100, Best Distance: 70
     Iteration 55/100, Best Distance: 70
     Iteration 56/100, Best Distance: 70
     Iteration 57/100, Best Distance: 70
     Iteration 58/100, Best Distance: 70
import numpy as np
from multiprocessing import Pool
import random
def two_opt(tour, distance_matrix):
    Performs the 2-opt optimization on the given tour.
    best tour = tour
    best_distance = calculate_distance(tour, distance_matrix)
    improved = True
    while improved:
        improved = False
        for i in range(1, len(best_tour) - 2):
            for j in range(i + 1, len(best_tour)):
                if j - i == 1: # Skip adjacent nodes
                new\_tour = best\_tour[:i] + best\_tour[i:j + 1][::-1] + best\_tour[j + 1:]
                new_distance = calculate_distance(new_tour, distance_matrix)
                if new_distance < best_distance:</pre>
                    best_tour = new_tour
                    best_distance = new_distance
                    improved = True
    return best_tour, best_distance
def calculate_distance(tour, distance_matrix):
    Calculates the total distance of the given tour.
    distance = 0
    for i in range(len(tour) - 1):
       distance += distance_matrix[tour[i]][tour[i + 1]]
    \label{linear_distance} \mbox{distance += distance\_matrix[tour[-1]][tour[0]]  \mbox{ \# Return to start}} \\
    return distance
class ParallelCellularOptimization:
    def __init__(self, distance_matrix, n_cells=4, n_iterations=100, alpha=0.5, migration_prob=0.2):
        self.distance_matrix = distance_matrix
        self.n_cells = n_cells
        self.n\_iterations = n\_iterations
        self.alpha = alpha
        self.migration_prob = migration_prob
    def _local_optimization(self, cell_solution):
        Applies local optimization using 2-opt for a single cell.
        return two_opt(cell_solution, self.distance_matrix)
    def _migrate(self, populations):
```

```
Handles migration between populations with the given probability.
        for i in range(len(populations) - 1):
            if random.random() < self.migration_prob:</pre>
                swap_idx = random.randint(0, len(populations[i]) - 1)
                populations[i][swap\_idx], \ populations[i+1][swap\_idx] = populations[i+1][swap\_idx], \ populations[i][swap\_idx] \\
        return populations
    def run(self):
        Runs the parallel cellular optimization.
        n_cities = len(self.distance_matrix)
        populations = [
           random.sample(range(n_cities), n_cities) for _ in range(self.n_cells)
        best_solution = None
       best distance = float('inf')
        for _ in range(self.n_iterations):
            # Parallel optimization
            with Pool(processes=self.n_cells) as pool:
                results = pool.map(self._local_optimization, populations)
            # Extract the best solution
            for solution, distance in results:
                if distance < best_distance:</pre>
                    best_solution = solution
                    best_distance = distance
            # Update populations for migration
            populations = [res[0] for res in results]
            populations = self. migrate(populations)
        return best_solution, best_distance
if __name__ == "__main__":
   # Example distance matrix (symmetric TSP)
    distance_matrix = np.random.randint(10, 100, size=(10, 10))
    np.fill diagonal(distance matrix, 0)
    # Ensure the matrix is symmetric
    distance_matrix = (distance_matrix + distance_matrix.T) // 2
    pco = ParallelCellularOptimization(distance_matrix, n_cells=4, n_iterations=100, alpha=0.5, migration_prob=0.2)
    best_solution, best_distance = pco.run()
    print(f"Best solution: {best_solution}")
    print(f"Best distance: {best_distance}")
    Best solution: [5, 2, 2, 2, 0, 3, 0, 4, 9, 9]
     Best distance: 207
import numpy as np
import random
class GeneticAlgorithm:
    def __init__(self, fitness_function, chromosome_length, population_size, mutation_rate, crossover_rate, generations):
        self.fitness_function = fitness_function
        self.chromosome_length = chromosome_length
        self.population_size = population_size
       self.mutation rate = mutation rate
        self.crossover_rate = crossover_rate
       self.generations = generations
    def _initialize_population(self):
        Randomly initialize the population with binary chromosomes.
        return np.random.randint(2, size=(self.population_size, self.chromosome_length))
    def _decode_chromosome(self, chromosome, lower_bound, upper_bound):
        Decode binary chromosome to a real number within a given range.
        decimal_value = int("".join(map(str, chromosome)), 2)
        max decimal = 2**self.chromosome length - 1
        return lower_bound + (upper_bound - lower_bound) * decimal_value / max_decimal
```

```
def _evaluate_fitness(self, population, lower_bound, upper_bound):
        Evaluate the fitness of the population.
        decoded_values = np.array(
            [self._decode_chromosome(chromosome, lower_bound, upper_bound) for chromosome in population]
        return np.array([self.fitness function(x) for x in decoded values])
    def _select_parents(self, population, fitness):
        Select parents using roulette wheel selection.
        # Adjust fitness to ensure non-negative probabilities
        adjusted_fitness = fitness - fitness.min() + 1e-6
        probabilities = adjusted_fitness / adjusted_fitness.sum()
        indices = np.random.choice(len(population), size=2, replace=False, p=probabilities)
        return population[indices[0]], population[indices[1]]
    def _crossover(self, parent1, parent2):
        Perform single-point crossover on two parents.
        if random.random() < self.crossover_rate:</pre>
            point = random.randint(1, self.chromosome_length - 1)
            child1 = np.concatenate((parent1[:point], parent2[point:]))
            child2 = np.concatenate((parent2[:point], parent1[point:]))
            return child1, child2
        return parent1.copy(), parent2.copy()
    def _mutate(self, chromosome):
        Mutate a chromosome by flipping bits with a given mutation rate.
        for i in range(len(chromosome)):
            if random.random() < self.mutation_rate:</pre>
                chromosome[i] = 1 - chromosome[i]
        return chromosome
    def optimize(self, lower_bound, upper_bound):
        Run the genetic algorithm to optimize the fitness function.
        population = self._initialize_population()
        best solution = None
        best_fitness = float("-inf")
        for generation in range(self.generations):
            fitness = self._evaluate_fitness(population, lower_bound, upper_bound)
            # Track the best solution
            max_fitness_idx = np.argmax(fitness)
            if fitness[max_fitness_idx] > best_fitness:
                best_fitness = fitness[max_fitness_idx]
                best_solution = population[max_fitness_idx]
            # Create the next generation
            new_population = []
            while len(new_population) < self.population_size:</pre>
                parent1, parent2 = self._select_parents(population, fitness)
                child1, child2 = self._crossover(parent1, parent2)
                new_population.append(self._mutate(child1))
                if len(new_population) < self.population_size:</pre>
                    new_population.append(self._mutate(child2))
            population = np.array(new_population)
            # Print progress
            print(f"Generation {generation + 1}: Best Fitness = {best_fitness:.4f}")
        best_decoded = self._decode_chromosome(best_solution, lower_bound, upper_bound)
        return best_decoded, best_fitness
if __name__ == "__main__":
   # Example fitness function: maximize f(x) = -x^2 + 5x + 10
    def fitness_function(x):
       return -x**2 + 5 * x + 10
    # GA parameters
    chromosome length = 16 # Precision of the solution
    population_size = 20
```

```
mutation_rate = 0.01
   crossover_rate = 0.8
   generations = 50
   lower bound = 0 # Lower bound of the search space
   upper_bound = 10 # Upper bound of the search space
   ga = GeneticAlgorithm(
       fitness_function, chromosome_length, population_size, mutation_rate, crossover_rate, generations
   best_solution, best_fitness = ga.optimize(lower_bound, upper_bound)
   print(f"\nBest solution: {best solution}")
   print(f"Best fitness: {best_fitness:.4f}")
→ Generation 1: Best Fitness = 16.1584
    Generation 2: Best Fitness = 16.1874
    Generation 3: Best Fitness = 16.1874
    Generation 4: Best Fitness = 16.1874
    Generation 5: Best Fitness = 16.1874
    Generation 6: Best Fitness = 16.2156
    Generation 7: Best Fitness = 16.2156
    Generation 8: Best Fitness = 16.2187
    Generation 9: Best Fitness = 16.2317
    Generation 10: Best Fitness = 16.2496
    Generation 11: Best Fitness = 16.2496
    Generation 12: Best Fitness = 16.2499
    Generation 13: Best Fitness = 16.2500
    Generation 14: Best Fitness = 16.2500
    Generation 15: Best Fitness = 16.2500
    Generation 16: Best Fitness = 16.2500
    Generation 17: Best Fitness = 16.2500
    Generation 18: Best Fitness = 16.2500
    Generation 19: Best Fitness = 16.2500
    Generation 20: Best Fitness = 16.2500
    Generation 21: Best Fitness = 16.2500
    Generation 22: Best Fitness = 16.2500
    Generation 23: Best Fitness = 16.2500
    Generation 24: Best Fitness = 16.2500
    Generation 25: Best Fitness = 16.2500
    Generation 26: Best Fitness = 16.2500
    Generation 27: Best Fitness = 16.2500
    Generation 28: Best Fitness = 16.2500
    Generation 29: Best Fitness = 16.2500
    Generation 30: Best Fitness = 16.2500
    Generation 31: Best Fitness = 16.2500
    Generation 32: Best Fitness = 16.2500
    Generation 33: Best Fitness = 16.2500
    Generation 34: Best Fitness = 16.2500
    Generation 35: Best Fitness = 16.2500
    Generation 36: Best Fitness = 16.2500
    Generation 37: Best Fitness = 16.2500
    Generation 38: Best Fitness = 16.2500
    Generation 39: Best Fitness = 16.2500
    Generation 40: Best Fitness = 16.2500
    Generation 41: Best Fitness = 16.2500
    Generation 42: Best Fitness = 16.2500
    Generation 43: Best Fitness = 16.2500
    Generation 44: Best Fitness = 16.2500
    Generation 45: Best Fitness = 16.2500
    Generation 46: Best Fitness = 16.2500
    Generation 47: Best Fitness = 16.2500
    Generation 48: Best Fitness = 16.2500
    Generation 49: Best Fitness = 16.2500
    Generation 50: Best Fitness = 16.2500
    Best solution: 2.5058365758754864
    Best fitness: 16.2500
```