```
#genetic algorithm
import numpy as np
import matplotlib.pyplot as plt
from sklearn.datasets import make_classification
from sklearn.model selection import train test split
from sklearn.metrics import accuracy_score
# Creating a sample dataset
X, y = make_classification(n_samples=500, n_features=10, n_informative=8, n_classes=2)
X_train, X_val, y_train, y_val = train_test_split(X, y, test_size=0.2)
# Neural Network Structure
input_size = X.shape[1]
hidden_size = 5
output_size = 1
# Helper functions for the Neural Network
def sigmoid(x):
    return 1 / (1 + np.exp(-x))
def forward_pass(X, weights1, weights2):
    hidden_input = np.dot(X, weights1)
    hidden_output = sigmoid(hidden_input)
    output_input = np.dot(hidden_output, weights2)
    output = sigmoid(output_input)
    return output
def compute_fitness(weights):
    predictions = forward\_pass(X\_train, weights['w1'], weights['w2'])
    predictions = (predictions > 0.5).astype(int)
    accuracy = accuracy_score(y_train, predictions)
    return accuracy
# Genetic Algorithm Parameters
population_size = 20
generations = 50
mutation_rate = 0.1
# Initialize Population
population = []
for _ in range(population_size):
    individual = {
        'w1': np.random.randn(input_size, hidden_size),
        'w2': np.random.randn(hidden_size, output_size)
    population.append(individual)
# Tracking performance
best_fitness_history = []
average_fitness_history = []
# Main Genetic Algorithm Loop
for generation in range(generations):
    # Evaluate Fitness of each Individual
    \verb|fitness_scores| = \verb|np.array|([compute_fitness(individual) | for individual | in population]||
    best_fitness = np.max(fitness_scores)
    average_fitness = np.mean(fitness_scores)
    best_fitness_history.append(best_fitness)
    average_fitness_history.append(average_fitness)
    # Selection: Select top half of the population
    sorted_indices = np.argsort(fitness_scores)[::-1]
    population = [population[i] for i in sorted_indices[:population_size//2]]
    # Crossover and Mutation
    new_population = []
    while len(new_population) < population_size:</pre>
        parents = np.random.choice(population, 2, replace=False)
        child = {
            'w1': (parents[0]['w1'] + parents[1]['w1']) / 2,
            'w2': (parents[0]['w2'] + parents[1]['w2']) / 2
        }
        # Mutation
        if np.random.rand() < mutation_rate:</pre>
            \label{eq:child['w1'] += np.random.randn(*child['w1'].shape) * 0.1} \\
            child['w2'] += np.random.randn(*child['w2'].shape) * 0.1
        new_population.append(child)
    population = new_population
```

```
print(f"Generation {generation+1}, Best Fitness: {best_fitness:.4f}")
# Evaluate the best individual on validation set
best_individual = population[np.argmax(fitness_scores)]
predictions = forward\_pass(X\_val, best\_individual['w1'], best\_individual['w2'])
predictions = (predictions > 0.5).astype(int)
final_accuracy = accuracy_score(y_val, predictions)
print(f"Final Accuracy on Validation Set: {final_accuracy:.4f}")
# Plotting the results
plt.figure(figsize=(10, 5))
plt.plot(best_fitness_history, label='Best Fitness')
plt.plot(average_fitness_history, label='Average Fitness')
plt.title('Fitness Over Generations')
plt.xlabel('Generation')
plt.ylabel('Fitness')
plt.legend()
plt.grid(True)
plt.show()
```

```
Generación i, pesc richess. 0.00//
    Generation 2, Best Fitness: 0.6775
    Generation 3, Best Fitness: 0.7350
    Generation 4, Best Fitness: 0.6975
    Generation 5, Best Fitness: 0.6925
    Generation 6, Best Fitness: 0.6950
    Generation 7, Best Fitness: 0.6975
    Generation 8, Best Fitness: 0.6975
    Generation 9, Best Fitness: 0.6975
    Generation 10, Best Fitness: 0.6975
    Generation 11, Best Fitness: 0.6975
    Generation 12, Best Fitness: 0.6950
    Generation 13, Best Fitness: 0.7000
    Generation 14, Best Fitness: 0.6975
    Generation 15, Best Fitness: 0.6975
    Generation 16, Best Fitness: 0.6975
    Generation 17, Best Fitness: 0.6975
    Generation 18, Best Fitness: 0.6975
    Generation 19, Best Fitness: 0.6975
    Generation 20, Best Fitness: 0.6975
    Generation 21, Best Fitness: 0.7000
    Generation 22, Best Fitness: 0.7000
    Generation 23, Best Fitness: 0.7050
    Generation 24, Best Fitness: 0.7000
    Generation 25, Best Fitness: 0.6975
    Generation 26, Best Fitness: 0.6975
Generation 27, Best Fitness: 0.6975
    Generation 28, Best Fitness: 0.7025
    Generation 29, Best Fitness: 0.7000
    Generation 30, Best Fitness: 0.7025
    Generation 31, Best Fitness: 0.7025
    Generation 32, Best Fitness: 0.7050
    Generation 33, Best Fitness: 0.7050
    Generation 34, Best Fitness: 0.7050
    Generation 35, Best Fitness: 0.7050
    Generation 36, Best Fitness: 0.7050
    Generation 37, Best Fitness: 0.7050
    Generation 38, Best Fitness: 0.7050
    Generation 39, Best Fitness: 0.7050
    Generation 40, Best Fitness: 0.7050
    Generation 41, Best Fitness: 0.7050
    Generation 42, Best Fitness: 0.7050
    Generation 43, Best Fitness: 0.7125
    Generation 44, Best Fitness: 0.7075
    Generation 45, Best Fitness: 0.7075
    Generation 46, Best Fitness: 0.7075
    Generation 47, Best Fitness: 0.7050
    Generation 48, Best Fitness: 0.7050
    Generation 49, Best Fitness: 0.7050
    Generation 50, Best Fitness: 0.7050
    Final Accuracy on Validation Set: 0.6200
```

0.70 0.65 0.60 0.55 0.50 0 10 20 30 40 50 Generation

```
# python implementation of particle swarm optimization (PSO)
# minimizing rastrigin and sphere function

import random
import math # cos() for Rastrigin
import copy # array-copying convenience
import sys # max float
```

```
#-----fitness functions-----
# rastrigin function
def fitness_rastrigin(position):
   fitnessVal = 0.0
    for i in range(len(position)):
       xi = position[i]
        fitnessVal += (xi * xi) - (10 * math.cos(2 * math.pi * xi)) + 10
    return fitnessVal
#sphere function
def fitness_sphere(position):
   fitnessVal = 0.0
    for i in range(len(position)):
       xi = position[i]
       fitnessVal += (xi*xi);
    return fitnessVal;
#particle class
class Particle:
   def __init__(self, fitness, dim, minx, maxx, seed):
        self.rnd = random.Random(seed)
        # initialize position of the particle with 0.0 value
        self.position = [0.0 for i in range(dim)]
       # initialize velocity of the particle with 0.0 value
        self.velocity = [0.0 for i in range(dim)]
        \# initialize best particle position of the particle with 0.0 value
        self.best_part_pos = [0.0 for i in range(dim)]
       # loop dim times to calculate random position and velocity
        # range of position and velocity is [minx, max]
        for i in range(dim):
            self.position[i] = ((maxx - minx) * self.rnd.random() + minx)
            self.velocity[i] = ((maxx - minx) * self.rnd.random() + minx)
        # compute fitness of particle
        self.fitness = fitness(self.position) # curr fitness
        # initialize best position and fitness of this particle
        self.best_part_pos = copy.copy(self.position)
        self.best_part_fitnessVal = self.fitness # best fitness
# particle swarm optimization function
def pso(fitness, max_iter, n, dim, minx, maxx):
   # hyper parameters
    w = 0.729 # inertia
   c1 = 1.49445 # cognitive (particle)
   c2 = 1.49445 \# social (swarm)
   rnd = random.Random(0)
   # create n random particles
   swarm = [Particle(fitness, dim, minx, maxx, i) for i in range(n)]
   \ensuremath{\text{\#}} compute the value of best_position and best_fitness in swarm
    best_swarm_pos = [0.0 for i in range(dim)]
    best_swarm_fitnessVal = sys.float_info.max # swarm best
    # computer best particle of swarm and it's fitness
    for i in range(n): # check each particle
        if swarm[i].fitness < best_swarm_fitnessVal:</pre>
            best_swarm_fitnessVal = swarm[i].fitness
            best_swarm_pos = copy.copy(swarm[i].position)
    # main loop of pso
    Iter = 0
    while Iter < max_iter:
        # after every 10 iterations
        # print iteration number and best fitness value so far
        if Iter % 10 == 0 and Iter > 1:
            print("Iter = " + str(Iter) + " best fitness = %.3f" % best_swarm_fitnessVal)
       for i in range(n): # process each particle
            # compute new velocity of curr particle
            for k in range(dim):
                r1 = rnd.random() # randomizations
```

```
r2 = rnd.random()
                swarm[i].velocity[k] = (
                                     (w * swarm[i].velocity[k]) +
                                     (c1 * r1 * (swarm[i].best_part_pos[k] - swarm[i].position[k])) +
                                     (c2 * r2 * (best_swarm_pos[k] -swarm[i].position[k]))
                # if velocity[k] is not in [minx, max]
                # then clip it
                if swarm[i].velocity[k] < minx:</pre>
                    swarm[i].velocity[k] = minx
                elif swarm[i].velocity[k] > maxx:
                   swarm[i].velocity[k] = maxx
            # compute new position using new velocity
            for k in range(dim):
                swarm[i].position[k] += swarm[i].velocity[k]
            # compute fitness of new position
            swarm[i].fitness = fitness(swarm[i].position)
            # is new position a new best for the particle?
            if swarm[i].fitness < swarm[i].best_part_fitnessVal:</pre>
                swarm[i].best_part_fitnessVal = swarm[i].fitness
                swarm[i].best_part_pos = copy.copy(swarm[i].position)
            # is new position a new best overall?
            if swarm[i].fitness < best_swarm_fitnessVal:</pre>
                best_swarm_fitnessVal = swarm[i].fitness
                best_swarm_pos = copy.copy(swarm[i].position)
        # for-each particle
       Iter += 1
    #end_while
    return best_swarm_pos
# end pso
#-----
# Driver code for rastrigin function
print("\nBegin particle swarm optimization on rastrigin function\n")
dim = 3
fitness = fitness_rastrigin
\label{eq:continuous}  \mbox{print("Goal is to minimize Rastrigin's function in " + str(dim) + " variables")} 
print("Function has known min = 0.0 at (", end="")
for i in range(dim-1):
    print("0, ", end="")
print("0)")
num_particles = 50
max iter = 100
print("Setting num_particles = " + str(num_particles))
print("Setting max_iter = " + str(max_iter))
print("\nStarting PSO algorithm\n")
best_position = pso(fitness, max_iter, num_particles, dim, -10.0, 10.0)
print("\nPSO completed\n")
print("\nBest solution found:")
print(["%.6f"%best_position[k] for k in range(dim)])
fitnessVal = fitness(best_position)
print("fitness of best solution = %.6f" % fitnessVal)
print("\\ \  nEnd\ particle\ swarm\ for\ rastrigin\ function\\ \  \  \  n")
print()
print()
# Driver code for Sphere function
print("\nBegin particle swarm optimization on sphere function\n")
dim = 3
```

```
fitness = fitness_sphere
print("Goal is to minimize sphere function in " + str(dim) + " variables")
print("Function has known min = 0.0 at (", end="")
for i in range(dim-1):
    print("0, ", end="")
print("0)")
num_particles = 50
max_iter = 100
print("Setting num_particles = " + str(num_particles))
print("Setting max_iter = " + str(max_iter))
print("\nStarting PSO algorithm\n")
best_position = pso(fitness, max_iter, num_particles, dim, -10.0, 10.0)
print("\nPSO completed\n")
print("\nBest solution found:")
print(["\%.6f"\%best\_position[k] \ for \ k \ in \ range(dim)])
fitnessVal = fitness(best_position)
print("fitness of best solution = %.6f" % fitnessVal)
print("\nEnd particle swarm for sphere function\n")
     Goal is to minimize Rastrigin's function in 3 variables
     Function has known min = 0.0 at (0, 0, 0)
     Setting num_particles = 50
     Setting max_iter = 100
     Starting PSO algorithm
     Iter = 10 best fitness = 8.463
     Iter = 20 best fitness = 4.792
     Iter = 30 best fitness = 2.223
     Iter = 40 best fitness = 0.251
     Iter = 50 best fitness = 0.251
     Iter = 60 best fitness = 0.061
     Iter = 70 best fitness = 0.007
     Iter = 80 best fitness = 0.005
     Iter = 90 best fitness = 0.000
     PSO completed
     Best solution found:
     ['0.000618', '0.000013', '0.000616']
     fitness of best solution = 0.000151
     End particle swarm for rastrigin function
     Begin particle swarm optimization on sphere function
     Goal is to minimize sphere function in 3 variables
     Function has known min = 0.0 at (0, 0, 0)
     Setting num_particles = 50
     Setting max_iter = 100
     Starting PSO algorithm
     Iter = 10 best fitness = 0.189
     Iter = 20 best fitness = 0.012
     Iter = 30 best fitness = 0.001
     Iter = 40 best fitness = 0.000
     Iter = 50 best fitness = 0.000
     Iter = 60 best fitness = 0.000
     Iter = 70 best fitness = 0.000
     Iter = 80 best fitness = 0.000
     Iter = 90 best fitness = 0.000
     PSO completed
     Best solution found:
     ['0.000004', '-0.000001', '0.000007']
fitness of best solution = 0.000000
     End particle swarm for sphere function
```

```
#ant colony
import numpy as np
import random
# Parameters
NUM_ANTS = 10
                      # Number of ants
NUM_CITIES = 5
                      # Number of cities
MAX_ITERATIONS = 100 # Number of iterations
ALPHA = 1.0
                      # Influence of pheromone
BETA = 2.0
                      # Influence of distance
EVAPORATION_RATE = 0.5 # Evaporation rate of pheromone
                      # Pheromone constant
INIT_PHEROMONE = 0.1 # Initial pheromone level
# Generate random distance matrix for TSP (symmetric graph)
def generate_distance_matrix(num_cities):
    distance_matrix = np.random.randint(1, 100, size=(num_cities, num_cities))
    np.fill_diagonal(distance_matrix, 0) \# Distance to itself is 0
    return (distance_matrix + distance_matrix.T) // 2 # Symmetric matrix
# Initialize pheromone matrix
def initialize_pheromones(num_cities, init_value):
    return np.full((num_cities, num_cities), init_value)
# Select next city based on probabilities
def select_next_city(current_city, visited, pheromones, distances):
    probabilities = []
    for city in range(len(distances)):
        if city not in visited:
            pheromone = pheromones[current_city][city] ** ALPHA
           heuristic = (1.0 / distances[current_city][city]) ** BETA
           probabilities.append(pheromone * heuristic)
        else:
           probabilities.append(0)
    probabilities = np.array(probabilities) / sum(probabilities)
    next_city = np.random.choice(range(len(distances)), p=probabilities)
    return next_city
# Update pheromones based on ant paths
def update pheromones(pheromones, paths, distances):
    pheromones *= (1 - EVAPORATION_RATE) # Evaporation
    for path in paths:
       path_length = calculate_path_length(path, distances)
        pheromone_deposit = Q / path_length
        for i in range(len(path) - 1):
           city_a, city_b = path[i], path[i + 1]
            pheromones[city_a][city_b] += pheromone_deposit
           pheromones[city_b][city_a] += pheromone_deposit # Symmetric
    return pheromones
# Calculate total length of a path
def calculate_path_length(path, distances):
   total_length = 0
    for i in range(len(path) - 1):
        total_length += distances[path[i]][path[i + 1]]
    total_length += distances[path[-1]][path[0]] # Return to start city
    return total_length
# Ant Colony Optimization
def ant_colony_optimization(num_cities, num_ants, max_iterations):
    distances = generate_distance_matrix(num_cities)
    pheromones = initialize_pheromones(num_cities, INIT_PHEROMONE)
    best path = None
    best_path_length = float('inf')
    for iteration in range(max_iterations):
        all_paths = []
        for ant in range(num_ants):
           visited = []
           current_city = random.randint(0, num_cities - 1)
           visited.append(current_city)
           while len(visited) < num cities:</pre>
                next_city = select_next_city(current_city, visited, pheromones, distances)
                visited.append(next_city)
                current_city = next_city
           visited.append(visited[0]) # Return to start city
           all_paths.append(visited)
```

```
path_length = calculate_path_length(visited, distances)
            if path_length < best_path_length:</pre>
               best_path = visited
               best_path_length = path_length
       # Update pheromones
        pheromones = update_pheromones(pheromones, all_paths, distances)
       print(f"Iteration {iteration + 1}: Best Path Length = {best_path_length}")
    print("\nBest Path:", best path)
    print("Best Path Length:", best_path_length)
# Run ACO for TSP
if __name__ == "__main__":
    ant_colony_optimization(NUM_CITIES, NUM_ANTS, MAX_ITERATIONS)
Iteration 2: Best Path Length = 216
     Iteration 3: Best Path Length = 216
     Iteration 4: Best Path Length = 216
     Iteration 5: Best Path Length = 216
     Iteration 6: Best Path Length = 216
     Iteration 7: Best Path Length = 216
     Iteration 8: Best Path Length = 216
     Iteration 9: Best Path Length = 216
     Iteration 10: Best Path Length = 216
     Iteration 11: Best Path Length = 216
     Iteration 12: Best Path Length = 216
     Iteration 13: Best Path Length = 216
     Iteration 14: Best Path Length = 216
     Iteration 15: Best Path Length = 216
     Iteration 16: Best Path Length = 216
     Iteration 17: Best Path Length = 216
     Iteration 18: Best Path Length = 216
     Iteration 19: Best Path Length = 216
     Iteration 20: Best Path Length = 216
     Iteration 21: Best Path Length = 216
     Iteration 22: Best Path Length = 216
     Iteration 23: Best Path Length = 216
     Iteration 24: Best Path Length = 216
     Iteration 25: Best Path Length = 216
     Iteration 26: Best Path Length = 216
     Iteration 27: Best Path Length = 216
     Iteration 28: Best Path Length = 216
     Iteration 29: Best Path Length = 216
     Iteration 30: Best Path Length = 216
     Iteration 31: Best Path Length = 216
     Iteration 32: Best Path Length = 216
     Iteration 33: Best Path Length = 216
     Iteration 34: Best Path Length = 216
     Iteration 35: Best Path Length = 216
     Iteration 36: Best Path Length = 216
     Iteration 37: Best Path Length = 216
     Iteration 38: Best Path Length = 216
     Iteration 39: Best Path Length = 216
     Iteration 40: Best Path Length = 216
     Iteration 41: Best Path Length = 216
     Iteration 42: Best Path Length = 216
     Iteration 43: Best Path Length = 216
     Iteration 44: Best Path Length = 216
     Iteration 45: Best Path Length = 216
     Iteration 46: Best Path Length = 216
     Iteration 47: Best Path Length = 216
     Iteration 48: Best Path Length = 216
     Iteration 49: Best Path Length = 216
     Iteration 50: Best Path Length = 216
     Iteration 51: Best Path Length = 216
     Iteration 52: Best Path Length = 216
     Iteration 53: Best Path Length = 216
     Iteration 54: Best Path Length = 216
     Iteration 55: Best Path Length = 216
     Iteration 56: Best Path Length = 216
     Iteration 57: Best Path Length = 216
     Iteration 58: Best Path Length = 216
import random
import networkx as nx
import numpy as np
import math
import datetime
import pandas as pd
class Cuckoo:
    def __init__(self, path, G, eps = 0.9):
```

```
self.path = path
              self.G = G
              self.nodes = list(G.nodes)
              self.eps = eps
              self.fitness = self.calculate_fitness()
       Function to Compute fitness value.
       def calculate_fitness(self):
              fitness = 0.0
              for i in range(1, len(self.path)):
                      total_distance = 0
                      curr_node = self.path[i-1]
                      next_node = self.path[i]
                      if self.G.has_edge(curr_node, next_node):
                            fitness += self.G[curr_node][next_node]['weight']
                             fitness += 0
              fitness = np.power(abs(fitness + self.eps), 2)
              return fitness
       def generate_new_path(self):
              This function generates a random solution (a random path) in the graph
              nodes = list(self.G.nodes)
              start = nodes[0]
              end = nodes[-1]
              samples = list(nx.all_simple_paths(self.G, start, end))
              for i in range(len(samples)):
                      if len(samples[i]) != len(nodes):
                             extra_nodes = [node for node in nodes if node not in samples[i]]
                             random.shuffle(extra_nodes)
                             samples[i] = samples[i] + extra_nodes
              sample_node = random.choice(samples)
              return sample_node
class CuckooSearch:
       def __init__(self, G, num_cuckoos, max_iterations, beta):
              self.G = G
              self.nodes = list(G.nodes)
              self.num_cuckoos = num_cuckoos
              self.max_iterations = max_iterations
              self.beta = beta
              self.cuckoos = [Cuckoo(random.sample(self.nodes, len(self.nodes)), self.G) for _ in range(self.num_cuckoos)]
              self.test_results = []
              self.test_cases = 0
       Function to buld new nests at new location and abandon old ones using Levi flights.
       def levy_flight(self):
              sigma = (math.gamma(1 + self.beta) * np.sin(np.pi * self.beta / 2) / (math.gamma((1 + self.beta) / 2) * self.beta * 2 ** ((self.beta) / 2) * (self.beta) * (self.beta) / 2) * (self.beta) * (self.beta) * (self.beta) / 2) * (self.beta) 
              u = np.random.normal(0, sigma, 1)
              v = np.random.normal(0, 1, 1)
              step = u / (abs(v) ** (1 / self.beta))
              return step
       def optimize(self):
               for i in range(self.max_iterations):
                      for j in range(self.num_cuckoos):
                             cuckoo = self.cuckoos[j]
                             step = self.levy_flight()
                             new_path = cuckoo.generate_new_path()
                             new_cuckoo = Cuckoo(new_path, self.G)
                             if new_cuckoo.fitness > cuckoo.fitness:
                                     self.cuckoos[j] = new_cuckoo
                                     self.test_cases+=1
                      self.cuckoos = sorted(self.cuckoos, key=lambda x: x.fitness, reverse=True)
                      best_path=self.cuckoos[0].path
                      best_fitness=self.cuckoos[0].fitness
                      self.test_results.append([i, best_fitness, self.test_cases])
               last_node = list(self.G.nodes)[-1]
              last_node_index = best_path.index(last_node) + 1
              return best_path[:last_node_index], best_fitness
```

```
if __name__ == "__main__":
    Example usage
    Gn = nx.DiGraph()
    #Add nodes to the graph
    for i in range(11):
        Gn.add_node(i)
    edges = [(0, 1, {\text{weight': 1}}), (1, 3, {\text{weight': 2}}), (1, 2, {\text{weight': 1}}), (2, 4, {\text{weight': 2}}),
             (3, 2,{'weight': 2}),(3, 4,{'weight': 1}),(3, 5,{'weight': 2}),(3, 7,{'weight': 4}),
             (4, 5,{'weight': 1}),(4, 6,{'weight': 2}),(5, 7,{'weight': 2}),(5, 8,{'weight': 3}),
             (6,\ 7,\{\text{'weight': 1}\}),(7,\ 9,\{\text{'weight': 2}\}),(8,\ 10,\{\text{'weight': 2}\}),(9,\ 10,\{\text{'weight': 1}\})]
    Gn.add_edges_from(edges)
    csa = CuckooSearch(Gn, num_cuckoos = 30, max_iterations=1000, beta=0.27)
    start = datetime.datetime.now()
    best_path, best_fitness = csa.optimize()
    end = datetime.datetime.now()
    csa_time = end - start
    csa_test_data = pd.DataFrame(csa.test_results, columns = ["iterations", "fitness_value", "test_cases"])
    print("Optimal path: ", best_path)
    print("Optimal path cost: ", best_fitness)
    print("CSA total Exec time => ", csa_time.total_seconds())
    csa_test_data.to_csv("csa_test_data_results.csv")
Optimal path: [0, 1, 3, 7, 9, 10]
Optimal path cost: 320.40999999999997
     CSA total Exec time => 12.944212
# Python implementation of Grey Wolf Optimization (GWO)
```

```
# Minimizing Rastrigin and Sphere function
import random
import math # cos() for Rastrigin
import copy # array-copying convenience
import sys
            # max float
#-----fitness functions-----
# Rastrigin function
def fitness_rastrigin(position):
    fitness_value = 0.0
    for i in range(len(position)):
       fitness_value += (xi * xi) - (10 * math.cos(2 * math.pi * xi)) + 10
    return fitness_value
# Sphere function
def fitness_sphere(position):
    fitness_value = 0.0
    for i in range(len(position)):
       xi = position[i]
       fitness_value += (xi * xi)
    return fitness_value
#-----
# Wolf class
class wolf:
    def __init__(self, fitness, dim, minx, maxx, seed):
       self.rnd = random.Random(seed)
       self.position = [0.0 for i in range(dim)]
       for i in range(dim):
           self.position[i] = ((maxx - minx) * self.rnd.random() + minx)
       self.fitness = fitness(self.position) # current fitness
# Grey Wolf Optimization (GWO)
def gwo(fitness, max_iter, n, dim, minx, maxx):
    rnd = random.Random(0)
    # Create n random wolves
    population = [wolf(fitness, dim, minx, maxx, i) for i in range(n)]
```

```
# On the basis of fitness values of wolves, sort the population in ascending order
    population = sorted(population, key=lambda temp: temp.fitness)
    # Best 3 solutions will be called as alpha, beta and gamma
    alpha_wolf, beta_wolf, gamma_wolf = copy.copy(population[:3])
    # Main loop of GWO
    Iter = 0
    while Iter < max_iter:</pre>
        # After every 10 iterations, print iteration number and best fitness value so far
        if Iter % 10 == 0 and Iter > 1:
            print("Iter = " + str(Iter) + " best fitness = %.3f" % alpha_wolf.fitness)
        # Linearly decreased from 2 to 0
        a = 2 * (1 - Iter / max_iter)
        # Updating each population member with the help of best three members
        for i in range(n):
            A1, A2, A3 = a * (2 * rnd.random() - 1), a * (2 * rnd.random() - 1), a * (2 * rnd.random() - 1)
            C1, C2, C3 = 2 * rnd.random(), 2 * rnd.random(), 2 * rnd.random()
            X1 = [0.0 \text{ for i in range(dim)}]
            X2 = [0.0 \text{ for i in range(dim)}]
            X3 = [0.0 \text{ for i in range(dim)}]
            Xnew = [0.0 for i in range(dim)]
            for j in range(dim):
                \label{eq:continuity} \textbf{X1[j]} = \texttt{alpha\_wolf.position[j]} - \texttt{A1} * \texttt{abs(C1} * \texttt{alpha\_wolf.position[j]} - \texttt{population[i].position[j])}
                X2[j] = beta_wolf.position[j] - A2 * abs(C2 * beta_wolf.position[j] - population[i].position[j])
                X3[j] = gamma\_wolf.position[j] - A3 * abs(C3 * gamma\_wolf.position[j] - population[i].position[j])
                Xnew[j] += X1[j] + X2[j] + X3[j]
            for j in range(dim):
                Xnew[j] /= 3.0
            # Fitness calculation of new solution
            fnew = fitness(Xnew)
            # Greedy selection
            if fnew < population[i].fitness:</pre>
                population[i].position = Xnew
                population[i].fitness = fnew
        # On the basis of fitness values of wolves, sort the population in ascending order
        population = sorted(population, key=lambda temp: temp.fitness)
        # Best 3 solutions will be called as alpha, beta, and gamma
        alpha_wolf, beta_wolf, gamma_wolf = copy.copy(population[:3])
        Iter += 1
    # End-while
    # Returning the best solution
    return alpha_wolf.position
# Driver code for Rastrigin function
print("\nBegin grey wolf optimization on Rastrigin function\n")
dim = 3
fitness = fitness_rastrigin
print("Goal is to minimize Rastrigin's function in " + str(dim) + " variables")
print("Function has known min = 0.0 at (", end="")
for i in range(dim-1):
    print("0, ", end="")
print("0)")
num particles = 50
max_iter = 100
print("Setting num_particles = " + str(num_particles))
print("Setting max_iter = " + str(max_iter))
print("\nStarting GWO algorithm\n")
best_position = gwo(fitness, max_iter, num_particles, dim, -10.0, 10.0)
print("\nGWO completed\n")
print("\nBest solution found:")
```

```
print(["%.6f" % best_position[k] for k in range(dim)])
err = fitness(best position)
print("Fitness of best solution = %.6f" % err)
print("\nEnd GWO for Rastrigin\n")
# Driver code for Sphere function
print("\nBegin grey wolf optimization on Sphere function\n")
dim = 3
fitness = fitness_sphere
print("Goal is to minimize Sphere function in " + str(dim) + " variables")
print("Function has known min = 0.0 at (", end="")
for i in range(dim-1):
   print("0, ", end="")
print("0)")
num particles = 50
max_iter = 100
print("Setting num_particles = " + str(num_particles))
print("Setting max_iter = " + str(max_iter))
print("\nStarting GWO algorithm\n")
best_position = gwo(fitness, max_iter, num_particles, dim, -10.0, 10.0)
print("\nGWO completed\n")
print("\nBest solution found:")
print(["%.6f" % best_position[k] for k in range(dim)])
err = fitness(best_position)
print("Fitness of best solution = %.6f" % err)
print("\nEnd GWO for Sphere\n")
\stackrel{	op}{\longrightarrow} Begin grey wolf optimization on Rastrigin function
     Goal is to minimize Rastrigin's function in 3 variables
     Function has known min = 0.0 at (0, 0, 0)
     Setting num_particles = 50
     Setting max_iter = 100
     Starting GWO algorithm
     Iter = 10 best fitness = 6.636
     Iter = 20 best fitness = 1.047
     Iter = 30 best fitness = 1.012
     Iter = 40 best fitness = 1.010
     Iter = 50 best fitness = 1.008
     Iter = 60 best fitness = 1.008
     Iter = 70 best fitness = 1.008
     Iter = 80 best fitness = 1.006
     Iter = 90 best fitness = 1.005
     GWO completed
     Best solution found:
     ['-0.004395', '0.995042', '0.005800']
     Fitness of best solution = 1.005465
     End GWO for Rastrigin
     Begin grey wolf optimization on Sphere function
     Goal is to minimize Sphere function in 3 variables
     Function has known min = 0.0 at (0, 0, 0)
     Setting num_particles = 50
     Setting max_iter = 100
     Starting GWO algorithm
     Iter = 10 best fitness = 0.000
     Iter = 20 best fitness = 0.000
     Iter = 30 best fitness = 0.000
     Iter = 40 best fitness = 0.000
     Iter = 50 best fitness = 0.000
     Iter = 60 best fitness = 0.000
     Iter = 70 best fitness = 0.000
     Iter = 80 best fitness = 0.000
     Iter = 90 best fitness = 0.000
     GWO completed
```

```
BEST SOLUTION TOURIG:

['0.000000', '0.000000', '-0.000000']

Fitness of best solution = 0.000000

End GWO for Sphere
```

```
import random
import multiprocessing
# Define the update rule function for each cell
# Here, a simple rule: New state = XOR of current cell and its two neighbors
def update_cell(index, grid, new_grid):
    # Handle the edge cells (leftmost and rightmost)
    left = grid[index - 1] if index > 0 else 0
    right = grid[index + 1] if index < len(grid) - 1 else 0
    # Rule: XOR of the current cell and its two neighbors
    new_grid[index] = left ^ grid[index] ^ right
# Parallel update function that processes all cells
def parallel_update(grid):
    # Create a new grid to store updated values
   new_grid = grid[:]
    # Create a pool of workers to update cells in parallel
    with multiprocessing.Pool(processes=multiprocessing.cpu\_count()) as pool:
        pool.starmap(update_cell, [(i, grid, new_grid) for i in range(len(grid))])
    return new_grid
# Function to run the automaton for a number of generations
def run_automaton(num_generations, grid_size):
    # Initialize the grid with random states (0 or 1)
    grid = [random.randint(0, 1) for _ in range(grid_size)]
    print("Initial Grid: ", grid)
    for generation in range(num_generations):
        print(f"Generation {generation + 1}: {grid}")
        grid = parallel_update(grid)
    return grid
          == '__main__':
if name
    grid_size = 20 # Size of the grid (number of cells)
    num_generations = 10 # Number of generations to run
    final_grid = run_automaton(num_generations, grid_size)
    print("Final Grid: ", final_grid)
Fr Initial Grid: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 1: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 2: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 3: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 4: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 5: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 6: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 7: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 8: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 9: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Generation 10: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
     Final Grid: [0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0]
import random
import string
```

```
import random
import string
import operator
import math

class GEP:
    def __init__(self, population_size, max_depth, max_generations, mutation_rate, crossover_rate):
        self.population_size = population_size
        self.max_depth = max_depth
        self.max_generations = max_generations
        self.mutation_rate = mutation_rate
        self.crossover_rate = crossover_rate
        self.operators = ['+', '-', '*', '/', '**']
        self.terminals = ['x', 'y', '1', '2', '3', '4', '5']
        self.population = self.initialize_population()

def initialize_population(self):
        population = []
```

```
_ iii i alige(seii.bobatatioli_size).
       individual = self.generate_individual()
       population.append(individual)
   return population
def generate_individual(self):
   individual = []
    depth = random.randint(1, self.max_depth)
   for _ in range(depth):
       individual.append(self.generate_expression())
    return individual
def generate_expression(self):
   expression = random.choice(self.operators + self.terminals)
   return expression
def fitness(self, individual, x_val, y_val):
   # Create a valid expression string
   expression = ''.join(individual)
   try:
       # Replace x and y with their values
       expression = expression.replace('x', str(x val)).replace('y', str(y val))
       # Evaluate the expression and return its absolute value as fitness (for simplicity)
        result = eval(expression)
       if math.isinf(result) or math.isnan(result):
            return float('inf') # Return high fitness for invalid results
       return abs(result)
   except:
       # In case of an error in evaluation (e.g., division by zero), return a high fitness value
       return float('inf')
def selection(self):
   fitness_values = [self.fitness(ind, 1, 1) for ind in self.population]
   total_fitness = sum(fitness_values)
   # Ensure that total fitness is non-zero and finite
   if total_fitness == 0 or not math.isfinite(total_fitness):
       return random.choice(self.population)
   # Normalize fitness values and select an individual based on fitness
   selection_probs = [fitness / total_fitness for fitness in fitness_values]
   return random.choices(self.population, selection_probs)[0]
def crossover(self, parent1, parent2):
   if random.random() < self.crossover_rate:</pre>
       # Ensure that we have valid crossover points by checking lengths
         crossover\_point = random.randint(1, \ min(len(parent1), \ len(parent2)) \ - \ 1) \ if \ min(len(parent1), \ len(parent2)) \ > \ 1 \ else \ 0 
       if crossover_point == 0:
           \# If crossover point is 0, return parents unchanged
            return parent1, parent2
       offspring1 = parent1[:crossover_point] + parent2[crossover_point:]
       offspring2 = parent2[:crossover_point] + parent1[crossover_point:]
        return offspring1, offspring2
   return parent1, parent2
def mutation(self, individual):
   if random.random() < self.mutation_rate:</pre>
       mutation_point = random.randint(0, len(individual) - 1)
       individual[mutation_point] = self.generate_expression()
    return individual
def evolve(self):
   best_fitness = float('inf')
   best solution = None
    for generation in range(self.max_generations):
        print(f"Generation {generation + 1}/{self.max_generations}")
       new population = []
        for _ in range(self.population_size // 2):
           # Selection
            parent1 = self.selection()
           parent2 = self.selection()
            offspring1, offspring2 = self.crossover(parent1, parent2)
            offspring1 = self.mutation(offspring1)
            offspring2 = self.mutation(offspring2)
            # Add offspring to the new population
            new_population.extend([offspring1, offspring2])
       # Replace the old population with the new population
```

```
self.population = new_population
            # Evaluate the population and find the best fitness
            current_best_fitness = float('inf')
            for individual in self.population:
                fitness_value = self.fitness(individual, 1, 1)
                if fitness_value < current_best_fitness:</pre>
                    current_best_fitness = fitness_value
                    best_solution = individual
            print(f"Best fitness so far: {current_best_fitness}")
            best_fitness = current_best_fitness
        return best_solution
# Initialize GEP with parameters and evolve the population
\verb|gep = GEP(population\_size=50, max\_depth=3, max\_generations=50, mutation\_rate=0.1, crossover\_rate=0.8)|
best_solution = gep.evolve()
print("Best solution found:")
print(best_solution)
→ Generation 1/50
     Best fitness so far: 1
     Generation 2/50
     Best fitness so far: 1
     Generation 3/50
     Best fitness so far: 4
     Generation 4/50
     Best fitness so far: 1
     Generation 5/50
```