

# **VISVESVARAYA TECHNOLOGICAL UNIVERSITY**

**“JnanaSangama”, Belgaum - 590014, Karnataka.**



## **LAB RECORD**

### **Bio Inspired Systems (23CS5BSBIS)**

*Submitted by*

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*in partial fulfillment for the award of the degree of*

**BACHELOR OF ENGINEERING  
*in*  
COMPUTER SCIENCE AND ENGINEERING**



**B.M.S. COLLEGE OF ENGINEERING  
(Autonomous Institution under VTU)  
BENGALURU-560019  
Aug-2025 to Dec-2025**

**B.M.S. College of Engineering,  
Bull Temple Road, Bangalore 560019**  
(Affiliated To Visvesvaraya Technological University, Belgaum)  
**Department of Computer Science and Engineering**



**CERTIFICATE**

This is to certify that the Lab work entitled “Bio Inspired Systems (23CS5BSBIS)” carried out by **Ruqaiyya Mahreen (1BM23CS351)**, who is a bonafide student of **B.M.S. College of Engineering**. It is in partial fulfillment for the award of **Bachelor of Engineering in Computer Science and Engineering** of the Visvesvaraya Technological University, Belgaum. The Lab report has been approved as it satisfies the academic requirements of the above mentioned subject and the work prescribed for the said degree.

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Github Link: <https://github.com/Ruqaiyya1BM23CS351/BISLab>

## Program 1

Implement a Genetic Algorithm to find the shortest possible route that visits a set of cities exactly once and returns to the starting city, optimizing the path using selection, crossover, and mutation operations.

Algorithm:

31/08/25 Week 1

classmate  
Date \_\_\_\_\_  
Page \_\_\_\_\_

**GENETIC ALGORITHM**

**Problem statement:** Travelling Salesman Problem  
Find the shortest path that visits each city exactly once and returns to the starting city.

**Fitness function:**

$$\text{fitness} = \frac{1}{\text{tot-dist}} \quad \text{since tot-dist has to be minimized.}$$

1) Initialize parameters:  
Population size: 50  
Generations: 500  
Crossover rate: 0.8  
Mutation rate: 0.2

2) Each individual is a random permutation of cities.

3) Selection -  
~~Roulette wheel~~ selection based on fitness

4) Crossover -  
Since chromosomes are permutations, we use ordered chromosomes.  
Select a slice from parent 1 and preserve order of remaining cities from parent 2.

5) Mutation:  
Swap mutation: swap position of 2 cities with a probability

6) Iteration:  
Evaluate → Selection → Crossover → Mutation → Next generation

Code:

```
import java.util.*;
```

```

public class GeneticTSP {

    // Parameters
    static final int POP_SIZE = 50;
    static final int GENS = 500;
    static final double CROSS_RATE = 0.8;
    static final double MUT_RATE = 0.2;

    static final int NUM_CITIES = 5;
    static final double[][] cities = {
        {0, 0}, {1, 5}, {5, 2}, {6, 6}, {8, 3}
    };

    static Random rand = new Random();

    // Individual representation
    static class Individual {
        int[] route;
        double fitness;

        Individual(int[] route) {
            this.route = route.clone();
            this.fitness = evaluate(route);
        }
    }

    // Euclidean distance
    static double distance(double[] c1, double[] c2) {
        return Math.sqrt(Math.pow(c1[0]-c2[0], 2) + Math.pow(c1[1]-c2[1], 2));
    }

    // Total distance of a route
    static double totalDistance(int[] route) {
        double dist = 0;
        for(int i = 0; i < NUM_CITIES; i++) {
            dist += distance(cities[route[i]], cities[route[(i+1)%NUM_CITIES]]);
        }
        return dist;
    }

    // Fitness function
    static double evaluate(int[] route) {
        return 1.0 / totalDistance(route);
    }
}

```

```

}

// Initialize population
static List<Individual> initPopulation() {
    List<Individual> population = new ArrayList<>();
    int[] base = new int[NUM_CITIES];
    for(int i=0;i<NUM_CITIES;i++) base[i]=i;
    for(int i=0;i<POP_SIZE;i++) {
        int[] route = base.clone();
        shuffle(route);
        population.add(new Individual(route));
    }
    return population;
}

// Shuffle array
static void shuffle(int[] array) {
    for(int i=array.length-1;i>0;i--) {
        int j = rand.nextInt(i+1);
        int temp = array[i];
        array[i] = array[j];
        array[j] = temp;
    }
}

// Tournament selection
static Individual select(List<Individual> population) {
    int tournamentSize = 5;
    Individual best = null;
    for(int i=0;i<tournamentSize;i++) {
        Individual ind = population.get(rand.nextInt(POP_SIZE));
        if(best==null || ind.fitness > best.fitness) best = ind;
    }
    return best;
}

// Ordered Crossover (OX)
static Individual[] crossover(Individual p1, Individual p2) {
    if(rand.nextDouble() > CROSS_RATE) {
        return new Individual[]{ new Individual(p1.route), new Individual(p2.route) };
    }
    int start = rand.nextInt(NUM_CITIES);
    int end = rand.nextInt(NUM_CITIES);
    if(start > end) { int tmp=start; start=end; end=tmp; }

```

```

        int[] c1 = new int[NUM_CITIES];
        Arrays.fill(c1, -1);
        int[] c2 = new int[NUM_CITIES];
        Arrays.fill(c2, -1);

        // Copy slice
        for(int i=start;i<end;i++) { c1[i]=p1.route[i]; c2[i]=p2.route[i]; }

        // Fill remaining
        fillRemaining(c1, p2.route, end);
        fillRemaining(c2, p1.route, end);

        return new Individual[]{ new Individual(c1), new Individual(c2) };
    }

    static void fillRemaining(int[] child, int[] parent, int start) {
        int idx = start;
        for(int i=0;i<NUM_CITIES;i++) {
            int city = parent[i];
            if(!contains(child, city)) {
                while(child[idx%NUM_CITIES]!=-1) idx++;
                child[idx%NUM_CITIES]=city;
            }
        }
    }

    static boolean contains(int[] arr, int val) {
        for(int v: arr) if(v==val) return true;
        return false;
    }

    // Swap Mutation
    static void mutate(Individual ind) {
        for(int i=0;i<NUM_CITIES;i++) {
            if(rand.nextDouble()<MUT_RATE) {
                int j = rand.nextInt(NUM_CITIES);
                int temp = ind.route[i];
                ind.route[i] = ind.route[j];
                ind.route[j] = temp;
            }
        }
        ind.fitness = evaluate(ind.route);
    }
}

```

```

public static void main(String[] args) {
    List<Individual> population = initPopulation();
    Individual best = null;

    for(int gen=0;gen<GENS;gen++) {
        List<Individual> newPop = new ArrayList<>();

        // Track best
        for(Individual ind: population) {
            if(best==null || ind.fitness>best.fitness) best=ind;
        }

        // Generate new population
        while(newPop.size()<POP_SIZE) {
            Individual p1 = select(population);
            Individual p2 = select(population);
            Individual[] offspring = crossover(p1, p2);
            mutate(offspring[0]);
            mutate(offspring[1]);
            newPop.add(offspring[0]);
            if(newPop.size()<POP_SIZE) newPop.add(offspring[1]);
        }
        population = newPop;
    }

    // Output best
    System.out.println("Best Route: " + Arrays.toString(best.route));
    System.out.println("Shortest Distance: " + totalDistance(best.route));
}
}

```

Output:

- Ruqaiyya@Ruqaiyya-mac ~ % cd "/Users/absk/ticTSP
   
Best Route: [4, 2, 0, 1, 3]
   
Shortest Distance: 22.35103276995244

## Program 2

Implement a Genetic Algorithm to solve the 0/1 Knapsack Problem — selecting a combination of items to maximize total value without exceeding the knapsack's weight capacity, using operations like selection, crossover, and mutation.

Algorithm:

12/9/25  
Week 2  
Gene Expression Algorithm

Problem statement - Knapsack (0/1)  
Each item has a weight and a value and a knapsack has a maximum weight capacity.  
Objective - select subset of items such that value is maximised without exceeding the capacity.

1) Fitness function:  
$$f(x) = \sum_{i=1}^n v_i x_i \quad x_i \in \{0, 1\}$$

2) Initialize the parameters -  
Population size: 4  
Number of genes: 4  
Mutation rate: 0.8  
Crossover rate: 0.1  
No of generations: 3

Item	Value	Weight
1	10	5
2	40	4
3	30	6
4	50	3

Capacity = 10

3) Each chromosome is a solution  
1010 (11 > 10)  
0101  
1110  
1001

4) Evaluate fitness  
Chromosome 1010 : Items {1, 3} → Value = 10 + 30 = 40  
Weight = 5 + 6 = 11 (exceeds)  
Fitness = 0

0101 : Items {2, 4} → Value = 40 + 50 = 90  
Weight = 4 + 3 = 7  
Fitness = 90

1110 : Value =  $10 + 40 + 30 = 80$   
 Weight =  $5 + 4 + 6 = 15$   
 Fitness = 20

1001 : Value = 60    Weight = 8    Fitness = 60  
 Best in generation 0 = 0101 (fitness 90)

5) Selection

choose parents based on fitness (0101 and 1001)

6) Crossover

One pt crossover between :

Parent 1: 0101

Parent 2: 1001

Cut after 2<sup>nd</sup> gene

Child 1 = 01/01 + 00/01 → 0101 { same as parents

Child 2 = 10/01 + 01/01 → 1001

7) Mutation

Suppose child 2 mutates at last gene

1001 → 1000

8) Gene expression

0101 → Items {2, 4} → Value = 90, Weight = 7

1000 → Items {1} → Value = 10, Weight = 5

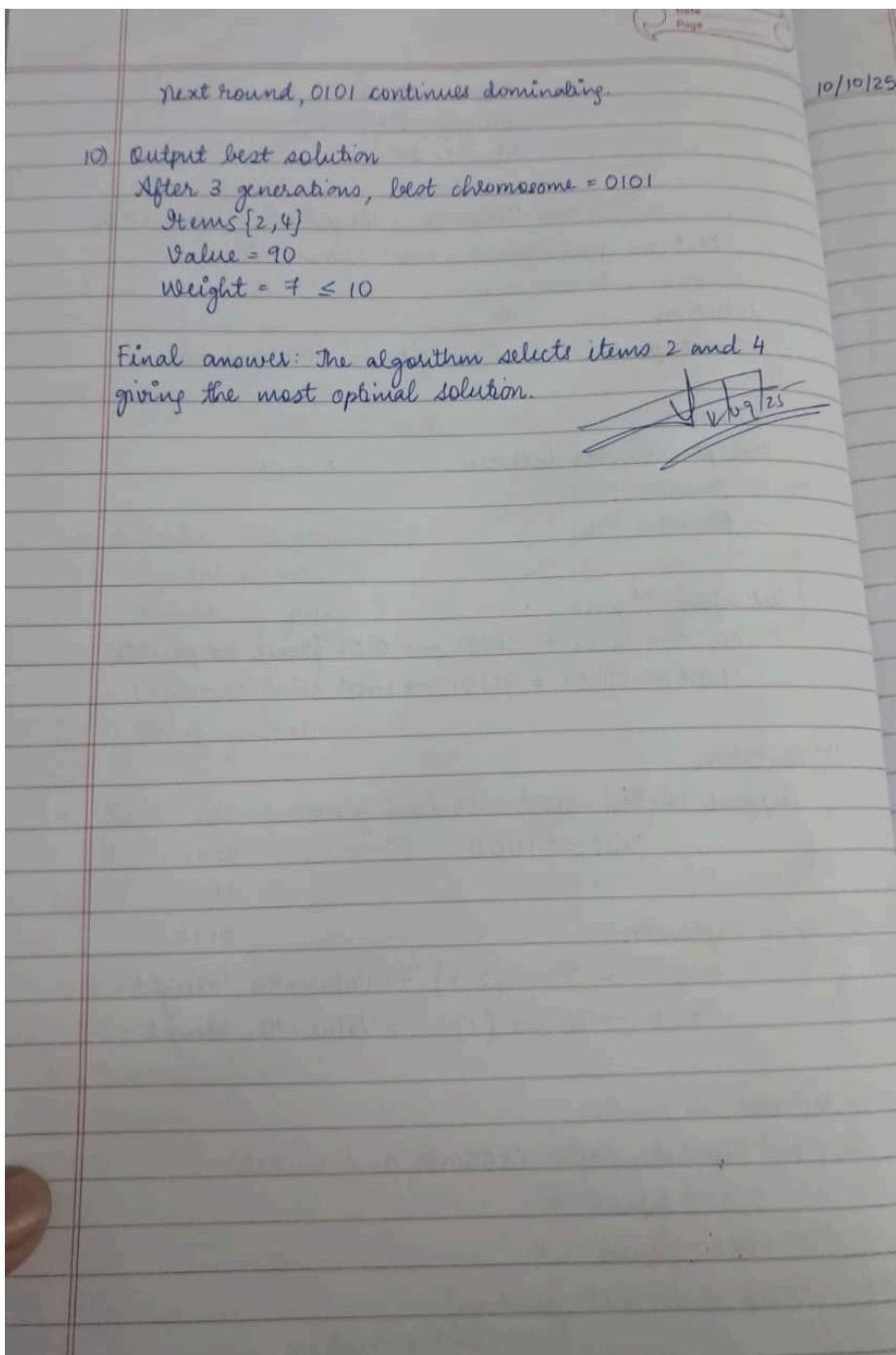
9) Iterate

New population (after crossover and mutation) :

0101 (fitness 90)

1000 (fitness 10)

Keep elites (best selection)  
from last generation



Code:

```
import random

# -----
# 1. Define the Problem
# -----
# Example items
values = [10, 40, 30, 50]    # values of the 4 items
weights = [5, 4, 6, 3]        # weights of the 4 items
capacity = 10                  # maximum knapsack weight
num_items = len(values)
```

```

# -----
# 2. Initialize Parameters
# -----
POP_SIZE = 10          # population size
GENERATIONS = 30        # number of generations
CROSSOVER_RATE = 0.8
MUTATION_RATE = 0.1

# -----
# 3. Initialize Population
# -----
def create_chromosome():
    """Generate a random binary chromosome."""
    return [random.randint(0, 1) for _ in range(num_items)]

def initialize_population():
    """Create an initial population."""
    return [create_chromosome() for _ in range(POP_SIZE)]

# -----
# 4. Evaluate Fitness
# -----
def fitness(chromosome):
    total_value = 0
    total_weight = 0
    for gene, v, w in zip(chromosome, values, weights):
        if gene == 1:
            total_value += v
            total_weight += w
    if total_weight > capacity:
        return 0 # penalty for exceeding capacity
    return total_value

# -----
# 5. Selection
# -----
def selection(population):
    """Tournament selection."""
    tournament_size = 3
    selected = random.sample(population, tournament_size)
    selected = sorted(selected, key=fitness, reverse=True)
    return selected[0]

# -----
# 6. Crossover
# -----
def crossover(parent1, parent2):
    if random.random() < CROSSOVER_RATE:
        point = random.randint(1, num_items - 1)
        child1 = parent1[:point] + parent2[point:]
        child2 = parent2[:point] + parent1[point:]
        return child1, child2
    return parent1[:], parent2[:]

# -----
# 7. Mutation

```

```

# -----
def mutate(chromosome):
    for i in range(len(chromosome)):
        if random.random() < MUTATION_RATE:
            chromosome[i] = 1 - chromosome[i] # flip bit
    return chromosome

# -----
# 8. Gene Expression (Phenotype Mapping)
# -----
def decode(chromosome):
    """Translate chromosome into selected items, total value, and total weight."""
    chosen_items = []
    total_value = 0
    total_weight = 0
    for i, gene in enumerate(chromosome):
        if gene == 1:
            chosen_items.append(i + 1)
            total_value += values[i]
            total_weight += weights[i]
    return chosen_items, total_value, total_weight

# -----
# 9. Iterate Through Generations
# -----
def gene_expression_algorithm():
    population = initialize_population()
    best_solution = None
    best_fitness = 0

    for generation in range(GENERATIONS):
        new_population = []

        # Evaluate all chromosomes
        population = sorted(population, key=fitness, reverse=True)
        if fitness(population[0]) > best_fitness:
            best_solution = population[0]
            best_fitness = fitness(population[0])

        # Print progress
        print(f"Generation {generation+1}: Best Fitness = {best_fitness}")

        # Elitism: carry the best two
        new_population.extend(population[:2])

        # Create new population
        while len(new_population) < POP_SIZE:
            parent1 = selection(population)
            parent2 = selection(population)
            child1, child2 = crossover(parent1, parent2)
            new_population.append(mutate(child1))
            if len(new_population) < POP_SIZE:
                new_population.append(mutate(child2))

        population = new_population

    return best_solution, best_fitness

```

```

# -----
# 10. Output Best Solution
# -----
best_chromosome, best_value = gene_expression_algorithm()
items, total_val, total_wt = decode(best_chromosome)

print("\nBest Chromosome:", best_chromosome)
print("Selected Items:", items)
print("Total Value:", total_val)
print("Total Weight:", total_wt)

```

Output:

```

Generation 1: Best Fitness = 70
Generation 2: Best Fitness = 90
Generation 3: Best Fitness = 90
Generation 4: Best Fitness = 90
Generation 5: Best Fitness = 90
Generation 6: Best Fitness = 90
Generation 7: Best Fitness = 90
Generation 8: Best Fitness = 90
Generation 9: Best Fitness = 90
Generation 10: Best Fitness = 90
Generation 11: Best Fitness = 90
Generation 12: Best Fitness = 90
Generation 13: Best Fitness = 90
Generation 14: Best Fitness = 90
Generation 15: Best Fitness = 90
Generation 16: Best Fitness = 90
Generation 17: Best Fitness = 90
Generation 18: Best Fitness = 90
Generation 19: Best Fitness = 90
Generation 20: Best Fitness = 90
Generation 21: Best Fitness = 90
Generation 22: Best Fitness = 90
Generation 23: Best Fitness = 90
Generation 24: Best Fitness = 90
Generation 25: Best Fitness = 90
Generation 26: Best Fitness = 90
Generation 27: Best Fitness = 90
Generation 28: Best Fitness = 90
Generation 29: Best Fitness = 90
Generation 30: Best Fitness = 90

Best Chromosome: [0, 1, 0, 1]
Selected Items: [2, 4]
Total Value: 90
Total Weight: 7

```

### Program 3

Implement the Particle Swarm Optimization algorithm to minimize a continuous mathematical function (e.g., the Sphere function) by simulating the collective behavior of particles that explore the search space to find the global optimum.

Algorithm:

10/10/23  
Particle Swarm Optimization

Initialization

- 1) Function : PSO( $S, W, C_1, C_2$ )
- 2) Input:
  - $S$  - Swarm size (no of particles)
  - $W$  - Inertia weight
  - $C_1$  - Cognitive (personal) constant
  - $C_2$  - Social (global) constant
- 3) Main loop:
  - for each particle  $i = 1$  to  $S$ :
    - initialize position ( $x_i$ )  $\in$  randomly within search space
    - initialize velocity ( $v_i$ ) randomly (often near zero)
    - (personal best position)  $PBest_i \leftarrow x_i$
    - $Fit(PBest_i) \leftarrow f(x_i)$
    - $GBest \leftarrow \text{ParticleWithBestFitness}(PBest)$

PSO Main loop:

Repeat until max iterations or convergence:

// Update best positions

for each particle  $i = 1$  to  $S$ :

- current fitness  $\leftarrow f(x_i)$
- if current fitness  $> Fit(PBest_i)$ :
  - $PBest_i \leftarrow x_i$
  - $Fit(PBest_i) \leftarrow \text{CurrentFitness}$
  - $GBest \leftarrow \text{ParticleWithBestFitness}(PBest)$

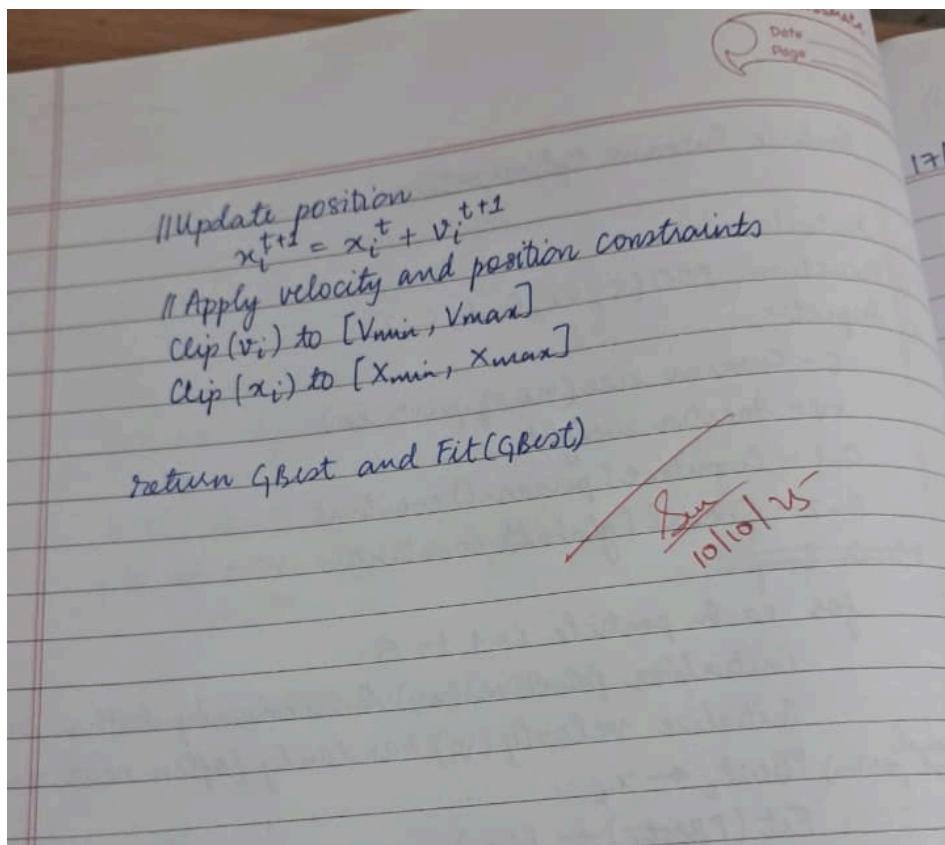
// Update velocity and position

for each particle  $i = 1$  to  $S$ :

- $random_1 \leftarrow \text{random}(0,1)$
- $random_2 \leftarrow \text{random}(0,1)$

// Update velocity

$$v_i^{t+1} = W \cdot v_i^t + C_1 \cdot random_1 \cdot (PBest_i - x_i^t) + C_2 \cdot random_2 \cdot (GBest - x_i^t)$$



Code:

```

import numpy as np
import random
import matplotlib.pyplot as plt
# Set seeds for reproducibility
random.seed(42)
np.random.seed(42)
# --- 1. Objective Function ---
def sphere_function(position):
    """
    The classic Sphere function ( $f(x) = \sum(x^2)$ ), used for minimization.
    The global minimum is  $f(x)=0$  at  $x=[0, 0, \dots, 0]$ .
    """
    # Ensures the input is treated as a NumPy array for vectorized operation
    return np.sum(position**2)

# --- 2. PSO Algorithm Implementation ---
def pso_optimizer(
    objective_func,
    num_particles=30,
    dimensions=2,
    search_range=(-10, 10),
    max_iterations=100,
    w=0.729,      # Inertia Weight (W)
    c1=1.4944,    # Cognitive Constant (C1)
    c2=1.4944    # Social Constant (C2)
):
    """
    Particle Swarm Optimization (PSO) algorithm for continuous optimization.

```

```

Args:
    objective_func (callable): The function to minimize.
    num_particles (int): Number of particles (S).
    dimensions (int): Dimensionality of the search space.
    search_range (tuple): (min, max) bounds for particle positions.
    max_iterations (int): Maximum number of generations.
    w (float): Inertia weight.
    c1 (float): Cognitive constant.
    c2 (float): Social constant.

"""
min_bound, max_bound = search_range

# 1. Initialize particle positions and velocities (x_i and v_i)
# Positions: [num_particles, dimensions]
positions = np.random.uniform(min_bound, max_bound, (num_particles, dimensions))
# Velocities: [num_particles, dimensions]
velocities = np.random.uniform(-1, 1, (num_particles, dimensions))
# 2. Initialize Personal Best (PBest_i)
pbest_positions = positions.copy()
pbest_scores = np.array([objective_func(p) for p in positions])
# 3. Initialize Global Best (GBest)
gbest_index = np.argmin(pbest_scores)
gbest_position = pbest_positions[gbest_index].copy()
gbest_score = pbest_scores[gbest_index]
history = [(gbest_score, gbest_position)]
print(f"Starting PSO for {dimensions} dimensions with {num_particles} particles...")
print(f"Initial GBest Score: {gbest_score:.4f}")
for iteration in range(max_iterations):
    # --- Phase 1: Update PBest Positions ---
    for i in range(num_particles):
        current_score = objective_func(positions[i])
        # Check if current position is better than particle's personal best
        if current_score < pbest_scores[i]:
            pbest_scores[i] = current_score
            pbest_positions[i] = positions[i].copy()

    # --- Update GBest (Global Best) ---
    # Find the overall best position among all PBest's
    current_gbest_index = np.argmin(pbest_scores)
    current_gbest_score = pbest_scores[current_gbest_index]

    # Update GBest only if a better PBest was found
    if current_gbest_score < gbest_score:
        gbest_score = current_gbest_score
        gbest_position = pbest_positions[current_gbest_index].copy()
    # Record history for convergence plot
    history.append((gbest_score, gbest_position.copy()))
    # --- Phase 2: Update Velocity and Position ---

    # Generate two sets of random numbers R1 and R2
    r1 = np.random.rand(num_particles, dimensions) # Random_1
    r2 = np.random.rand(num_particles, dimensions) # Random_2
    # 1. Inertia component: W * v_i^t
    inertia_comp = w * velocities
    # 2. Cognitive component (PBest influence): C1 * r1 * (PBest_i - x_i^t)
    cognitive_comp = c1 * r1 * (pbest_positions - positions)
    # 3. Social component (GBest influence): C2 * r2 * (GBest - x_i^t)

```

```

        # NumPy handles broadcasting of the 1D gbest_position to the 2D positions matrix
        social_comp = c2 * r2 * (gbest_position - positions)
        # Update velocity: v_i^{t+1} = Inertia + Cognitive + Social
        velocities = inertia_comp + cognitive_comp + social_comp
        # Update position: x_i^{t+1} = x_i^t + v_i^{t+1}
        positions = positions + velocities
        # Apply position constraints (clipping to search space)
        positions = np.clip(positions, min_bound, max_bound)
        print(f"Iteration {iteration+1}/{max_iterations}: GBest Score =
{gbest_score:.4e}")
    return gbest_position, gbest_score, history
# --- 3. Example Usage and Visualization ---
def run_pso_example():
    # Set up PSO parameters
    search_range = (-5.12, 5.12) # Common range for Sphere function
    max_iter = 100
    # Run the PSO solver
    best_position, best_score, history = pso_optimizer(
        objective_func=sphere_function,
        num_particles=30,
        dimensions=2, # Using 2D for simple visualization
        search_range=search_range,
        max_iterations=max_iter
    )
    print("\n--- Results ---")
    print(f"Objective Function: Sphere Function")
    print(f"Best Position Found: {best_position}")
    print(f"Minimum Score (Fitness): {best_score:.6e}")
    # --- Visualization ---
    # Extract scores for convergence plot
    scores = [item[0] for item in history]

    plt.figure(figsize=(10, 5))

    # Plot 1: Convergence History
    plt.subplot(1, 2, 1)
    plt.plot(range(len(scores)), scores, color='darkorange', linewidth=2)
    plt.title('PSO Convergence History')
    plt.xlabel('Iteration')
    plt.ylabel('Global Best Score (Log Scale)', color='darkorange')
    plt.yscale('log') # Use log scale for better visualization of minimization
    plt.grid(True, which="both", ls="--")
    # Plot 2: Particle movement (only for 2D problems)
    if history and len(history[0][1]) == 2:
        plt.subplot(1, 2, 2)

        # Create a contour plot of the objective function
        x = np.linspace(search_range[0], search_range[1], 100)
        y = np.linspace(search_range[0], search_range[1], 100)
        X, Y = np.meshgrid(x, y)

        # Calculate Z values for the Sphere function across the grid
        Z = np.array([[sphere_function(np.array([X[i, j], Y[i, j]])) for j in range(100)] for i in range(100)])

        plt.contourf(X, Y, Z, levels=50, cmap='viridis')
        plt.colorbar(label='Function Value')

```

```
# Plot the final best position
plt.plot(best_position[0], best_position[1], 'r*', markersize=15, label='Final
GBest')
plt.title('2D Search Space Visualization')
plt.xlabel('Dimension 1 (X)')
plt.ylabel('Dimension 2 (Y)')
plt.legend()
plt.tight_layout()
plt.show()

# Execute the example
if __name__ == '__main__':
    run_pso_example()
```

Output:

Starting PSO for 2 dimensions with 50 particles....

```

Totalt Objekt Score: 1.795
Dimension 1/100: Objekt Score = 1.795e+00
Dimension 2/100: Objekt Score = 1.830e-01
Dimension 3/100: Objekt Score = 1.747e-01
Dimension 4/100: Objekt Score = 1.747e-01
Dimension 5/100: Objekt Score = 1.747e-01
Dimension 6/100: Objekt Score = 1.747e-01
Dimension 7/100: Objekt Score = 1.747e-01
Dimension 8/100: Objekt Score = 1.747e-01
Dimension 9/100: Objekt Score = 1.747e-01
Dimension 10/100: Objekt Score = 1.830e-01
Dimension 11/100: Objekt Score = 2.052e-01
Dimension 12/100: Objekt Score = 1.703e-01
Dimension 13/100: Objekt Score = 1.703e-01
Dimension 14/100: Objekt Score = 1.953e-01
Dimension 15/100: Objekt Score = 1.953e-01
Dimension 16/100: Objekt Score = 1.953e-01
Dimension 17/100: Objekt Score = 1.953e-01
Dimension 18/100: Objekt Score = 2.052e-01
Dimension 19/100: Objekt Score = 2.052e-01
Dimension 20/100: Objekt Score = 1.771e-01
Dimension 21/100: Objekt Score = 1.771e-01
Dimension 22/100: Objekt Score = 1.771e-01
Dimension 23/100: Objekt Score = 1.800e-01
Dimension 24/100: Objekt Score = 1.800e-01
Dimension 25/100: Objekt Score = 1.800e-01
Dimension 26/100: Objekt Score = 1.800e-01
Dimension 27/100: Objekt Score = 1.800e-01
Dimension 28/100: Objekt Score = 1.800e-01
Dimension 29/100: Objekt Score = 1.800e-01
Dimension 30/100: Objekt Score = 1.800e-01
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Dimension 38/100: Objekt Score = 1.800e-01
Dimension 39/100: Objekt Score = 1.800e-01
Dimension 40/100: Objekt Score = 1.800e-01
Dimension 41/100: Objekt Score = 1.800e-01
Dimension 42/100: Objekt Score = 1.800e-01
Dimension 43/100: Objekt Score = 1.800e-01
Dimension 44/100: Objekt Score = 7.149e-02
Dimension 45/100: Objekt Score = 7.149e-02
Dimension 46/100: Objekt Score = 7.149e-02
Dimension 47/100: Objekt Score = 4.732e-02
Dimension 48/100: Objekt Score = 4.732e-02
Dimension 49/100: Objekt Score = 4.732e-02
Dimension 50/100: Objekt Score = 4.732e-02
Dimension 51/100: Objekt Score = 4.732e-02
Dimension 52/100: Objekt Score = 4.732e-02
Dimension 53/100: Objekt Score = 4.732e-02
Dimension 54/100: Objekt Score = 4.732e-02
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Dimension 56/100: Objekt Score = 4.732e-02
Dimension 57/100: Objekt Score = 4.732e-02
Dimension 58/100: Objekt Score = 4.732e-02
Dimension 59/100: Objekt Score = 4.732e-02
Dimension 60/100: Objekt Score = 4.732e-02
Dimension 61/100: Objekt Score = 4.732e-02
Dimension 62/100: Objekt Score = 4.732e-02
Dimension 63/100: Objekt Score = 4.732e-02
Dimension 64/100: Objekt Score = 4.732e-02
Dimension 65/100: Objekt Score = 4.732e-02
Dimension 66/100: Objekt Score = 4.732e-02
Dimension 67/100: Objekt Score = 4.732e-02
Dimension 68/100: Objekt Score = 4.732e-02
Dimension 69/100: Objekt Score = 4.732e-02
Dimension 70/100: Objekt Score = 4.732e-02
Dimension 71/100: Objekt Score = 4.732e-02
Dimension 72/100: Objekt Score = 2.795e-02
Dimension 73/100: Objekt Score = 2.795e-02
Dimension 74/100: Objekt Score = 2.795e-02
Dimension 75/100: Objekt Score = 2.795e-02
Dimension 76/100: Objekt Score = 2.795e-02
Dimension 77/100: Objekt Score = 2.381e-02
Dimension 78/100: Objekt Score = 1.469e-02
Dimension 79/100: Objekt Score = 1.469e-02
Dimension 80/100: Objekt Score = 1.469e-02
Dimension 81/100: Objekt Score = 1.469e-02
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Dimension 95/100: Objekt Score = 1.469e-02
Dimension 96/100: Objekt Score = 1.469e-02
Dimension 97/100: Objekt Score = 1.469e-02
Dimension 98/100: Objekt Score = 1.469e-02
Dimension 99/100: Objekt Score = 1.469e-02
Dimension 100/100: Objekt Score = 1.777e-13

```

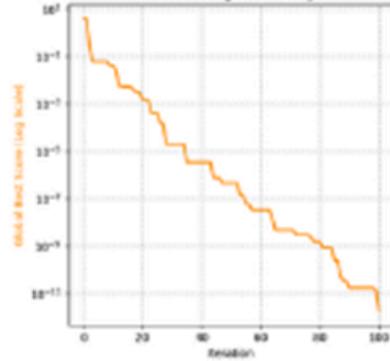
— AreaObj

Objective Function: Sphere Function

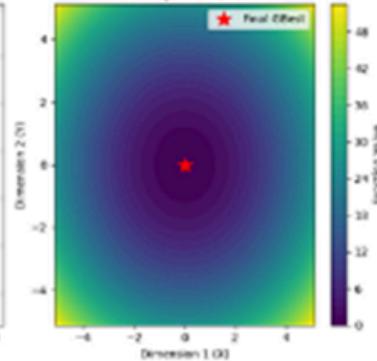
Best Position Found: [0.0352644e-07 0.2076034e-07]

Minimum Score (Fitness): 1.77754e-13

PSO Convergence History



2D Search Space Visualization



## Program 4

Implement the Ant Colony Optimization algorithm to find the shortest possible route that visits all cities exactly once and returns to the starting city (Traveling Salesman Problem).

Algorithm:

10/10/25 Week 3  
Ant Colony Optimization -  
For TSP-

Initialization:

- 1) Function: ACO-TSP(M, N, T-0,  $\alpha$ ,  $\beta$ ,  $\rho$ )
- 2) Input:
  - M: No of ants (particles)
  - N: No of cities (nodes)
  - T-0: Initial pheromone value on all edges
  - $\alpha$ : Pheromone influence parameter
  - $\beta$ : Heuristic (distance) influence parameter
  - $\rho$ : Pheromone evaporation rate ( $0 < \rho < 1$ )
- 3) Initialize pheromone ( $T_{ij}$ ) on all edges  $(i, j)$  to  $T-0$
- 4) Best Tour  $\leftarrow$  NULL
- 5) Best length  $\leftarrow$  Infinity

ACO Main loop:  
Repeat until max iterations:  
//Tour construction  
for each ant  $k = 1$  to  $M$ :  
    Ant $_k$ .Tour  $\leftarrow$  startAtRandomCity()  
    while Ant $_k$  has not visited all cities:  
        current city  $\leftarrow$  last city in (Ant $_k$ .tour)  
        select next city  $j$  from unvisited cities  $J$  based on  
        probability  $P_{ij}^k$   
$$P_{ij}^k = \frac{[T_{ij}]^\alpha \cdot [n_{ij}]^\beta}{\sum_{i \in J} [T_{ij}]^\alpha [n_{ij}]^\beta}$$
  
        //  $T_{ij}$  - pheromone trail  
        //  $n_{ij} = \frac{1}{\text{Distance } (i, j)}$   
        Ant $_k$ .tour  $\leftarrow$  Append(j)  
    Ant $_k$ .length  $\leftarrow$  calculateTotDist(Ant $_k$ .tour)

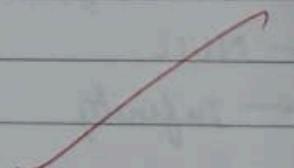
```

if ant_k.length < bestLength:
    bestLength ← Ant_k.length
    bestTour ← Ant_k.tour

// Pheromone update
for all edges (i, j):
    Tij ← (1 - ρ) · Tij
    ΔT ← 1 / bestLength
    for all edges (i, j) in bestTour:
        Tij ← Tij + ΔT

return bestTour and bestLength

```



Code:

```

import numpy as np
import random
import matplotlib.pyplot as plt

# --- 1. Utility Functions ---
def create_distance_matrix(cities):
    """Calculates the Euclidean distance matrix between all pairs of cities."""
    num_cities = len(cities)
    dist_matrix = np.zeros((num_cities, num_cities))
    for i in range(num_cities):
        for j in range(i + 1, num_cities):
            # Calculate Euclidean distance
            distance = np.sqrt((cities[i][0] - cities[j][0])**2 + (cities[i][1] - cities[j][1])**2)
            dist_matrix[i, j] = dist_matrix[j, i] = distance
    return dist_matrix

```

```

def calculate_tour_length(tour, dist_matrix):
    """Calculates the total length of a given tour (sequence of city indices)."""
    length = 0
    num_cities = len(tour)
    for i in range(num_cities):
        # Add distance from current city to next city in the tour
        city_a = tour[i]
        city_b = tour[(i + 1) % num_cities] # Wrap around to the start city
        length += dist_matrix[city_a, city_b]
    return length

# --- 2. ACO Algorithm Implementation ---
def aco_tsp_solver(cities, num_ants=10, max_iterations=100, alpha=1.0, beta=5.0, rho=0.5,
initial_pheromone=1.0):
    """
    Ant Colony Optimization (ACO) algorithm for the Traveling Salesman Problem (TSP).
    Args:
        cities (list of tuples): List of (x, y) coordinates for each city.
        num_ants (int): Number of artificial ants (M).
        max_iterations (int): Maximum number of generations to run.
        alpha (float): Influence of the pheromone trail ( $\tau$ ).
        beta (float): Influence of the heuristic information ( $\eta$ ,  $1/distance$ ).
        rho (float): Pheromone evaporation rate.
        initial_pheromone (float): Initial pheromone value ( $\tau_0$ ).
    """
    num_cities = len(cities)
    dist_matrix = create_distance_matrix(cities)

    # Heuristic matrix ( $\eta_{ij} = 1 / distance_{ij}$ )
    eta_matrix = 1.0 / (dist_matrix + np.finfo(float).eps)
    np.fill_diagonal(eta_matrix, 0)

    # Initialize pheromone matrix ( $\tau_{ij}$ )
    pheromone_matrix = np.full((num_cities, num_cities), initial_pheromone)

    # Initialize best tour found so far
    best_tour = None
    best_length = float('inf')
    history = []

    print(f"Starting ACO for {num_cities} cities with {num_ants} ants...")

    for iteration in range(max_iterations):
        all_tours = []
        all_lengths = []

        # --- Phase 1: Tour Construction ---
        for ant in range(num_ants):
            start_city = random.randint(0, num_cities - 1)
            tour = [start_city]
            visited = {start_city}

            for _ in range(num_cities - 1):
                current_city = tour[-1]
                unvisited_cities = [c for c in range(num_cities) if c not in visited]

```

```

        if not unvisited_cities:
            break # Should not happen if TSP is solvable

        # Calculate probabilities P_ij
        probabilities = []
        denominator = 0.0

        for next_city in unvisited_cities:
            tau = pheromone_matrix[current_city, next_city] ** alpha
            eta = eta_matrix[current_city, next_city] ** beta
            numerator = tau * eta
            probabilities.append((next_city, numerator))
            denominator += numerator

        if denominator == 0:
            # Fallback to random choice if all probabilities are zero (rare)
            next_city = random.choice(unvisited_cities)
        else:
            # Select next city based on roulette wheel selection (weighted
            probability)
            prob_values = [p[1] / denominator for p in probabilities]
            next_city = random.choices(
                [p[0] for p in probabilities],
                weights=prob_values,
                k=1
            )[0]

            tour.append(next_city)
            visited.add(next_city)

        tour_length = calculate_tour_length(tour, dist_matrix)
        all_tours.append(tour)
        all_lengths.append(tour_length)

        # Update personal best (used for global best update below)
        if tour_length < best_length:
            best_length = tour_length
            best_tour = tour

        history.append(best_length)

    # --- Phase 2: Pheromone Update ---
    # 1. Evaporation: tau_ij = (1 - rho) * tau_ij
    pheromone_matrix = (1 - rho) * pheromone_matrix

    # 2. Deposition: The best ant deposits pheromone (Ant System variant)
    if best_tour is not None:
        # Pheromone deposit (Delta_tau = 1 / BestLength)
        delta_tau = 1.0 / best_length
        # Deposit pheromone along the best tour
        for i in range(num_cities):
            city_a = best_tour[i]
            city_b = best_tour[(i + 1) % num_cities]
            pheromone_matrix[city_a, city_b] += delta_tau
            pheromone_matrix[city_b, city_a] += delta_tau # TSP graph is symmetric

    print(f"Iteration {iteration+1}/{max_iterations}: Best Length =

```

```

{best_length:.2f}")

    return best_tour, best_length, history

# --- 3. Example Usage and Visualization ---
def run_aco_example():
    # Define a simple set of 10 cities (x, y coordinates)
    random.seed(42)  # for reproducibility
    np.random.seed(42)
    cities = [(random.uniform(0, 10), random.uniform(0, 10)) for _ in range(10)]

    # Run the ACO solver
    best_tour, best_length, history = aco_tsp_solver(
        cities,
        num_ants=20,
        max_iterations=50,
        alpha=1.0,
        beta=5.0,
        rho=0.1
    )

    print("\n--- Results ---")
    print(f"Cities: {cities}")
    print(f"Best Tour (City Indices): {best_tour}")
    print(f"Best Tour Length: {best_length:.4f}")

    # --- Visualization ---
    if best_tour:
        # Prepare coordinates for plotting the best tour path
        x_coords = [cities[i][0] for i in best_tour]
        y_coords = [cities[i][1] for i in best_tour]
        # Close the loop for visualization
        x_coords.append(x_coords[0])
        y_coords.append(y_coords[0])

        plt.figure(figsize=(10, 5))

        # Plot 1: Tour Path
        plt.subplot(1, 2, 1)
        plt.plot(x_coords, y_coords, 'o-', color='blue', markerfacecolor='red',
        markersize=8)
        # Label cities with their index
        for i, (x, y) in enumerate(cities):
            plt.text(x + 0.1, y + 0.1, str(i), fontsize=9)
        plt.title(f'ACO Best TSP Tour (Length: {best_length:.2f})')
        plt.xlabel('X Coordinate')
        plt.ylabel('Y Coordinate')
        plt.grid(True)

        # Plot 2: Convergence History
        plt.subplot(1, 2, 2)
        plt.plot(history, color='green', linewidth=2)
        plt.title('ACO Convergence')
        plt.xlabel('Iteration')
        plt.ylabel('Best Tour Length')
        plt.grid(True)

```

```
plt.tight_layout()  
plt.show()  
  
# Execute the example  
if __name__ == '__main__':  
    run_aco_example()
```

Output:

```

Starting ACO for 10 cities with 20 ants...
Iteration 1/50: Best Length = 27.17
Iteration 2/50: Best Length = 26.41
Iteration 3/50: Best Length = 26.41
Iteration 4/50: Best Length = 26.41
Iteration 5/50: Best Length = 26.41
Iteration 6/50: Best Length = 26.41
Iteration 7/50: Best Length = 26.41
Iteration 8/50: Best Length = 26.41
Iteration 9/50: Best Length = 26.41
Iteration 10/50: Best Length = 26.41
Iteration 11/50: Best Length = 26.41
Iteration 12/50: Best Length = 26.41
Iteration 13/50: Best Length = 26.41
Iteration 14/50: Best Length = 26.41
Iteration 15/50: Best Length = 26.41
Iteration 16/50: Best Length = 26.41
Iteration 17/50: Best Length = 26.41
Iteration 18/50: Best Length = 26.41
Iteration 19/50: Best Length = 26.41
Iteration 20/50: Best Length = 26.41
Iteration 21/50: Best Length = 26.41
Iteration 22/50: Best Length = 26.41
Iteration 23/50: Best Length = 26.41
Iteration 24/50: Best Length = 26.41
Iteration 25/50: Best Length = 26.41
Iteration 26/50: Best Length = 26.41
Iteration 27/50: Best Length = 26.41
Iteration 28/50: Best Length = 26.41
Iteration 29/50: Best Length = 26.41
Iteration 30/50: Best Length = 26.41
Iteration 31/50: Best Length = 26.41
Iteration 32/50: Best Length = 26.41
Iteration 33/50: Best Length = 26.41
Iteration 34/50: Best Length = 26.41
Iteration 35/50: Best Length = 26.41
Iteration 36/50: Best Length = 26.41
Iteration 37/50: Best Length = 26.41
Iteration 38/50: Best Length = 26.41
Iteration 39/50: Best Length = 26.41
Iteration 40/50: Best Length = 26.41
Iteration 41/50: Best Length = 26.41
Iteration 42/50: Best Length = 26.41
Iteration 43/50: Best Length = 26.41
Iteration 44/50: Best Length = 26.41
Iteration 45/50: Best Length = 26.41
Iteration 46/50: Best Length = 26.41
Iteration 47/50: Best Length = 26.41
Iteration 48/50: Best Length = 26.41
Iteration 49/50: Best Length = 26.41
Iteration 50/50: Best Length = 26.41

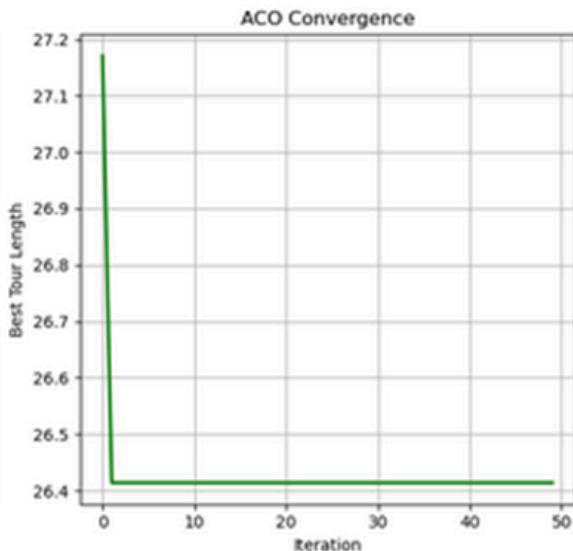
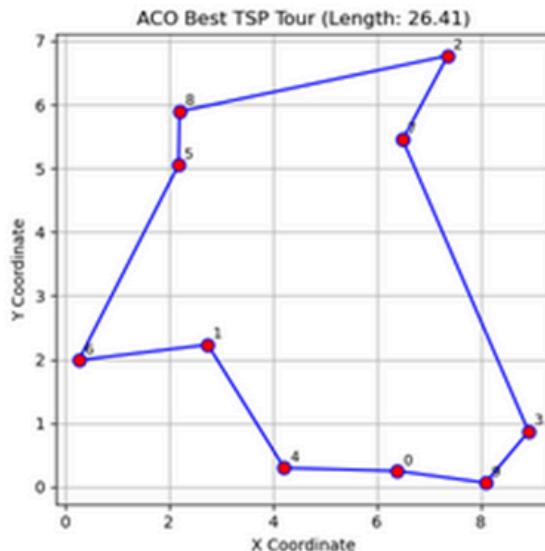
```

--- Results ---

```

Cities: [(6.394267084578837, 0.25810755222666936), (2.7582931836011926, 2.2321073814882277), (7.364712141648124, 6.766904874229113), (8.921795677048454, 8.8693883262041615), (4.2192181968527045, 0.297972194388078344), (2.1863797488360336, 5.053552881033624), (0.26535960683863625, 1.988376506866485), (6.408844377795232, 5.449414886632166), (2.284406226486067, 5.892656838759888), (8.004384566778266, 0.06408759678061617)]
Best Tour (City Indices): [7, 2, 8, 5, 6, 1, 4, 0, 9, 3]
Best Tour Length: 26.4139

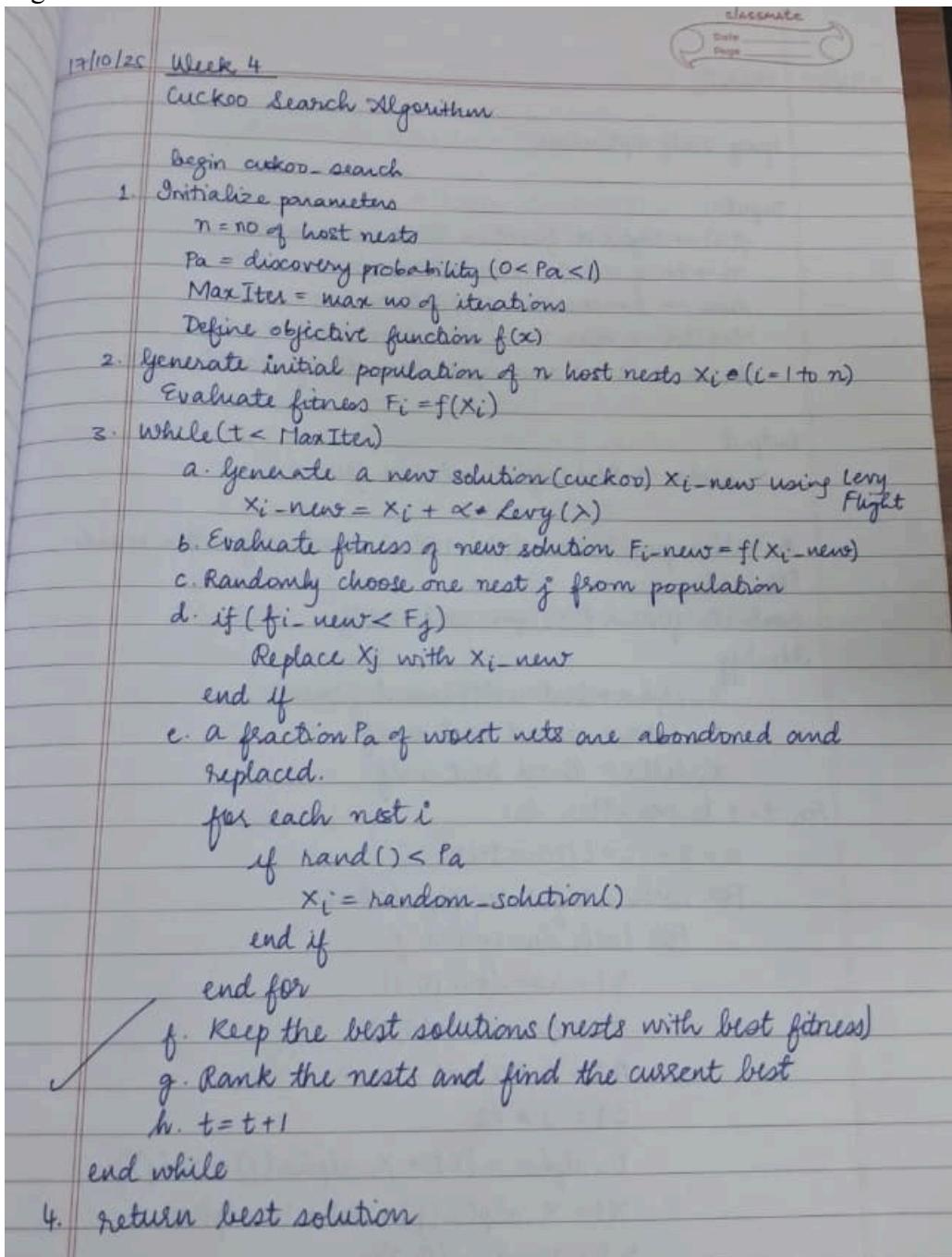
```



## Program 5

Use the Cuckoo Search metaheuristic to minimize a given objective function (e.g., Sphere function) by iteratively improving candidate solutions through Lévy flights and random nest replacement.

Algorithm:



17/10/25 Week 4  
Cuckoo Search Algorithm

classmate  
Date \_\_\_\_\_  
Page \_\_\_\_\_

```
begin cuckoo-search
1. Initialize parameters
    n = no of host nests
    Pa = discovery probability (0 < Pa < 1)
    MaxIter = max no of iterations
    Define objective function f(x)
2. Generate initial population of n host nests  $x_i \in (i=1 \text{ to } n)$ 
    Evaluate fitness  $F_i = f(x_i)$ 
3. While ( $t < \text{MaxIter}$ )
    a. Generate a new solution (cuckoo)  $x_i\text{-new}$  using Lévy Flight
         $x_i\text{-new} = x_i + \alpha * \text{Levy}(\lambda)$ 
    b. Evaluate fitness of new solution  $F_i\text{-new} = f(x_i\text{-new})$ 
    c. Randomly choose one nest  $j$  from population
    d. if ( $f_i\text{-new} < F_j$ )
        Replace  $x_j$  with  $x_i\text{-new}$ 
    end if
    e. a fraction  $Pa$  of worst nests are abandoned and replaced.
    for each nest  $i$ 
        if rand() <  $Pa$ 
             $x_i = \text{random\_solution}()$ 
        end if
    end for
    f. Keep the best solutions (nests with best fitness)
    g. Rank the nests and find the current best
    h.  $t = t + 1$ 
end while
4. return best solution
```

Code:

```
import numpy as np
import math

# Objective Function (example: minimize Sphere function)
def objective_function(x):
```

```

    return np.sum(x**2)

# Lévy flight function
def levy_flight(Lambda):
    # Using math module instead of np.math
    sigma = (math.gamma(1 + Lambda) * math.sin(math.pi * Lambda / 2) /
              (math.gamma((1 + Lambda) / 2) * Lambda * 2 ** ((Lambda - 1) / 2))) ** (1 / Lambda)
    u = np.random.randn() * sigma
    v = np.random.randn()
    step = u / abs(v) ** (1 / Lambda)
    return step

# Cuckoo Search Algorithm
def cuckoo_search(obj_func, dim=2, n=15, pa=0.25, alpha=0.01, max_iter=100):
    # Initialize nests randomly
    nests = np.random.uniform(-5, 5, size=(n, dim))
    fitness = np.array([obj_func(x) for x in nests])

    # Find the current best nest
    best_idx = np.argmin(fitness)
    best = nests[best_idx].copy()

    for t in range(max_iter):
        for i in range(n):
            # Generate a new solution via Lévy flight
            step_size = alpha * levy_flight(1.5)
            new_nest = nests[i] + step_size * (nests[i] - best)
            new_nest = np.clip(new_nest, -5, 5)

            # Evaluate fitness
            f_new = obj_func(new_nest)

            # Replace if the new solution is better
            if f_new < fitness[i]:
                nests[i] = new_nest
                fitness[i] = f_new

        # Abandon some nests (with probability pa)
        for i in range(n):
            if np.random.rand() < pa:
                nests[i] = np.random.uniform(-5, 5, dim)
                fitness[i] = obj_func(nests[i])

        # Update the best solution
        best_idx = np.argmin(fitness)
        if fitness[best_idx] < obj_func(best):
            best = nests[best_idx].copy()

        # Print progress
        print(f"Iteration {t+1}: Best fitness = {obj_func(best):.6f}")

    return best, obj_func(best)

# Run the algorithm
best_solution, best_value = cuckoo_search(objective_function, dim=2, n=20, max_iter=50)
print("\nBest Solution:", best_solution)
print("Best Fitness Value:", best_value)

```

**Output:**

```
Iteration 1: Best fitness = 0.108432
Iteration 2: Best fitness = 0.108432
Iteration 3: Best fitness = 0.108432
Iteration 4: Best fitness = 0.108432
Iteration 5: Best fitness = 0.108432
Iteration 6: Best fitness = 0.108432
Iteration 7: Best fitness = 0.108432
Iteration 8: Best fitness = 0.108432
Iteration 9: Best fitness = 0.108432
Iteration 10: Best fitness = 0.108432
Iteration 11: Best fitness = 0.108432
Iteration 12: Best fitness = 0.108432
Iteration 13: Best fitness = 0.108432
Iteration 14: Best fitness = 0.108432
Iteration 15: Best fitness = 0.108432
Iteration 16: Best fitness = 0.108432
Iteration 17: Best fitness = 0.108432
Iteration 18: Best fitness = 0.108432
Iteration 19: Best fitness = 0.108432
Iteration 20: Best fitness = 0.108432
Iteration 21: Best fitness = 0.108432
Iteration 22: Best fitness = 0.108432
Iteration 23: Best fitness = 0.108432
Iteration 24: Best fitness = 0.108432
Iteration 25: Best fitness = 0.108432
Iteration 26: Best fitness = 0.108432
Iteration 27: Best fitness = 0.108432
Iteration 28: Best fitness = 0.108432
Iteration 29: Best fitness = 0.108432
Iteration 30: Best fitness = 0.108432
Iteration 31: Best fitness = 0.108432
Iteration 32: Best fitness = 0.108432
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Iteration 35: Best fitness = 0.108432
Iteration 36: Best fitness = 0.108432
Iteration 37: Best fitness = 0.108432
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Iteration 40: Best fitness = 0.108432
Iteration 41: Best fitness = 0.108432
Iteration 42: Best fitness = 0.108432
Iteration 43: Best fitness = 0.108432
Iteration 44: Best fitness = 0.108432
Iteration 45: Best fitness = 0.108432
Iteration 46: Best fitness = 0.108432
Iteration 47: Best fitness = 0.108432
Iteration 48: Best fitness = 0.108432
Iteration 49: Best fitness = 0.108432
Iteration 50: Best fitness = 0.108432
```

Best Solution: [0.27812108 0.17629828]  
Best Fitness Value: 0.10843242171891111

## Program 6

Implement the Grey Wolf Optimizer algorithm to minimize a mathematical objective function (e.g.,  $f(x) = x^2 - 4x + 4$ ) by simulating the social hierarchy and hunting behavior of grey wolves to find the global optimum.

Algorithm:

5/11/25 week 5

### Grey Wolf Optimizer

Input:

- $f(x)$  → Objective function to be minimized
- $n$  → no. of wolves (population size)
- $dim$  → Dimension of problem
- $MaxIter$  → max no. of iterations
- $lb, ub$  → lower and upper bounds of search space

Output:

- $x_{\text{alpha}}$  → Best (alpha) solution found

Initialize population  $x_i$  ( $i=1$  to  $n$ ) randomly within bounds  $[lb, ub]$

Evaluate fitness  $f(x_i)$  for each wolf

Identify :

- $x_{\text{alpha}} = \text{best wolf (lowest fitness)}$
- $x_{\text{beta}} = \text{second best wolf}$
- $x_{\delta} = \text{third best wolf}$

For  $t = 1$  to  $MaxIter$  do:

$a = 2 - (2 * t / MaxIter)$

For each wolf  $i$  in population:

For each dimension  $j$ :

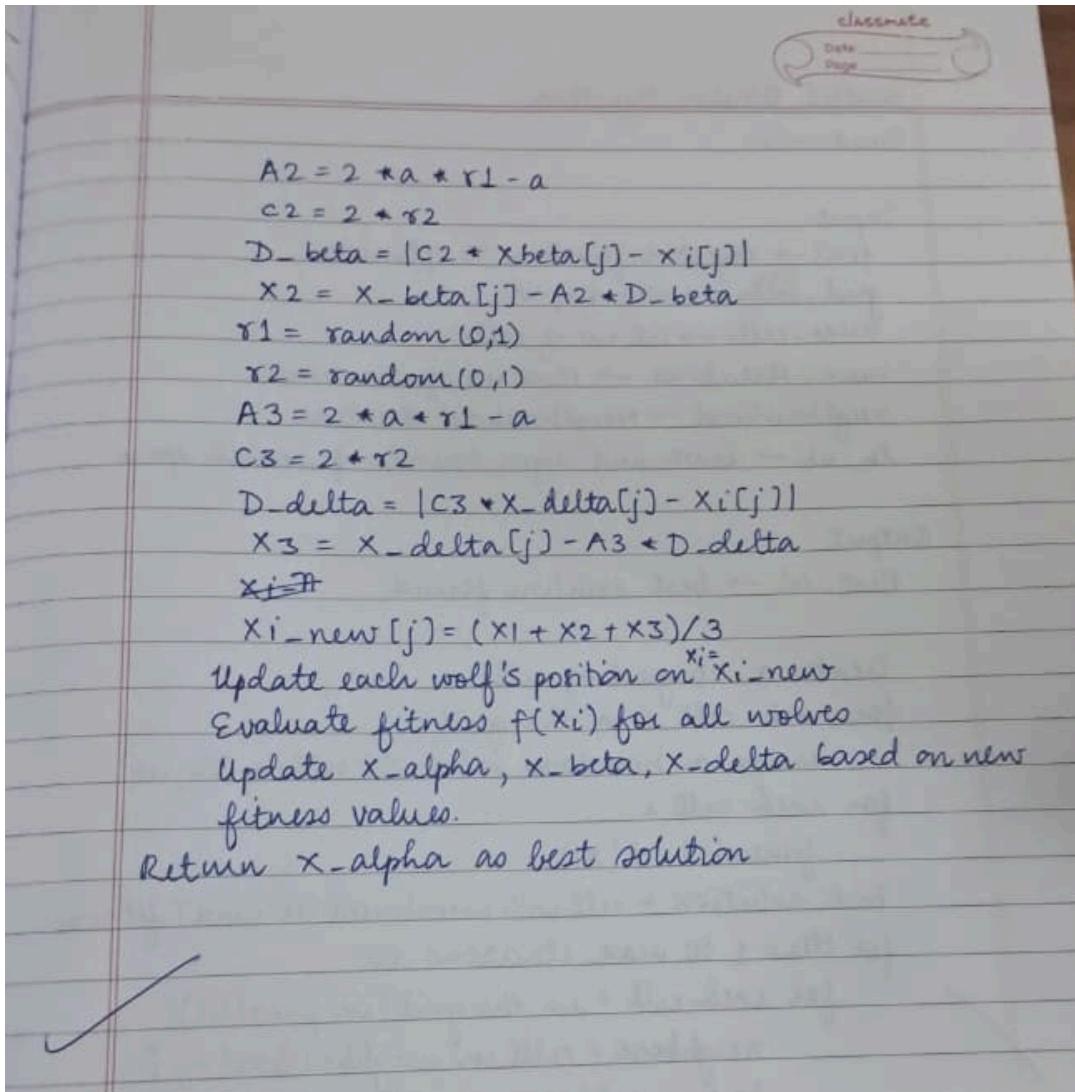
- $r_1 = \text{random}(0, 1)$
- $r_2 = \text{random}(0, 1)$
- $A_1 = 2 * a * r_1 - a$
- $C_1 = 2 * r_2$

$D_{\text{alpha}} = [C_1 * x_{\text{alpha}}[j] - x_i[j]]$

$X_1 = x_{\text{alpha}}[j] - A_1 + D_{\text{alpha}}$

$r_1 = \text{random}(0, 1)$

$r_2 = \text{random}(0, 1)$



Code:

```

import numpy as np

# Objective function (example: minimize f(x) = x^2 - 4x + 4)
def objective_function(x):
    return x[0]**2 - 4*x[0] + 4 # single-variable function

# Grey Wolf Optimizer
def GWO(obj_func, dim, n_wolves, max_iter, lb, ub):
    # Initialize positions of wolves randomly within [lb, ub]
    wolves = np.random.uniform(lb, ub, (n_wolves, dim))

    # Initialize alpha, beta, delta positions and fitness
    alpha_pos = np.zeros(dim)
    beta_pos = np.zeros(dim)
    delta_pos = np.zeros(dim)
    alpha_score = float("inf")
    beta_score = float("inf")
    delta_score = float("inf")
  
```

```

# Main optimization loop
for t in range(max_iter):
    for i in range(n_wolves):
        # Clip wolves to the boundary
        wolves[i] = np.clip(wolves[i], lb, ub)

        # Calculate fitness
        fitness = obj_func(wolves[i])

        # Update alpha, beta, delta
        if fitness < alpha_score:
            alpha_score, alpha_pos = fitness, wolves[i].copy()
        elif fitness < beta_score:
            beta_score, beta_pos = fitness, wolves[i].copy()
        elif fitness < delta_score:
            delta_score, delta_pos = fitness, wolves[i].copy()

    # Linearly decreasing coefficient 'a' from 2 to 0
    a = 2 - t * (2 / max_iter)

    # Update position of each wolf
    for i in range(n_wolves):
        for j in range(dim):
            r1, r2 = np.random.rand(), np.random.rand()
            A1 = 2 * a * r1 - a
            C1 = 2 * r2
            D_alpha = abs(C1 * alpha_pos[j] - wolves[i][j])
            X1 = alpha_pos[j] - A1 * D_alpha

            r1, r2 = np.random.rand(), np.random.rand()
            A2 = 2 * a * r1 - a
            C2 = 2 * r2
            D_beta = abs(C2 * beta_pos[j] - wolves[i][j])
            X2 = beta_pos[j] - A2 * D_beta

            r1, r2 = np.random.rand(), np.random.rand()
            A3 = 2 * a * r1 - a
            C3 = 2 * r2
            D_delta = abs(C3 * delta_pos[j] - wolves[i][j])
            X3 = delta_pos[j] - A3 * D_delta

            # Average of three best positions
            wolves[i][j] = (X1 + X2 + X3) / 3

        # Optional: Print progress
        print(f"Iteration {t+1}/{max_iter} → Best fitness: {alpha_score:.6f}")

    return alpha_pos, alpha_score

# Example usage
best_position, best_score = GWO(
    obj_func=objective_function,
    dim=1,                  # single-variable function
    n_wolves=10,             # number of wolves
    max_iter=50,              # number of iterations
    lb=-10,                 # lower bound
    ub=10                   # upper bound
)

```

```
)  
  
print("\nBest Solution Found:")  
print("x =", best_position)  
print("Fitness =", best_score)
```

Output:

```
Iteration 1/50 → Best fitness: 0.041188
Iteration 2/50 → Best fitness: 0.041188
Iteration 3/50 → Best fitness: 0.041188
Iteration 4/50 → Best fitness: 0.041188
Iteration 5/50 → Best fitness: 0.004941
Iteration 6/50 → Best fitness: 0.002758
Iteration 7/50 → Best fitness: 0.002758
Iteration 8/50 → Best fitness: 0.002758
Iteration 9/50 → Best fitness: 0.001457
Iteration 10/50 → Best fitness: 0.001457
Iteration 11/50 → Best fitness: 0.000078
Iteration 12/50 → Best fitness: 0.000078
Iteration 13/50 → Best fitness: 0.000040
Iteration 14/50 → Best fitness: 0.000040
Iteration 15/50 → Best fitness: 0.000040
Iteration 16/50 → Best fitness: 0.000040
Iteration 17/50 → Best fitness: 0.000040
Iteration 18/50 → Best fitness: 0.000040
Iteration 19/50 → Best fitness: 0.000040
Iteration 20/50 → Best fitness: 0.000040
Iteration 21/50 → Best fitness: 0.000040
Iteration 22/50 → Best fitness: 0.000040
Iteration 23/50 → Best fitness: 0.000040
Iteration 24/50 → Best fitness: 0.000040
Iteration 25/50 → Best fitness: 0.000007
Iteration 26/50 → Best fitness: 0.000007
Iteration 27/50 → Best fitness: 0.000007
Iteration 28/50 → Best fitness: 0.000007
Iteration 29/50 → Best fitness: 0.000007
Iteration 30/50 → Best fitness: 0.000007
Iteration 31/50 → Best fitness: 0.000007
Iteration 32/50 → Best fitness: 0.000007
Iteration 33/50 → Best fitness: 0.000007
Iteration 34/50 → Best fitness: 0.000007
Iteration 35/50 → Best fitness: 0.000005
Iteration 36/50 → Best fitness: 0.000005
Iteration 37/50 → Best fitness: 0.000005
Iteration 38/50 → Best fitness: 0.000005
Iteration 39/50 → Best fitness: 0.000005
Iteration 40/50 → Best fitness: 0.000005
Iteration 41/50 → Best fitness: 0.000005
Iteration 42/50 → Best fitness: 0.000002
Iteration 43/50 → Best fitness: 0.000001
Iteration 44/50 → Best fitness: 0.000001
Iteration 45/50 → Best fitness: 0.000001
Iteration 46/50 → Best fitness: 0.000001
Iteration 47/50 → Best fitness: 0.000001
Iteration 48/50 → Best fitness: 0.000001
Iteration 49/50 → Best fitness: 0.000001
Iteration 50/50 → Best fitness: 0.000001

Best Solution Found:
x = [2.00108665]
Fitness = 1.1808108792976668e-06
```

## Program 7

Implement the Parallel Cellular Algorithm to minimize a mathematical objective function (e.g.,  $f(x) = x^2 - 4x + 4$ ) by modeling a population of solutions arranged in a 2D grid, where each cell evolves based on the best solution in its local neighborhood to gradually approach the global optimum.

Algorithm:

Parallel Cellular Algorithm  
Pseudocode:

**Input:**  
 $f(x)$  → objective function to optimize  
grid\_size → size of grid  
num\_cells → tot no of cells  
max\_iterations → Max no of iterations  
neighbourhood → Neighbourhood structure  
 $lb, ub$  → lower and upper bounds of search space

**Output:**  
 $best\_sol \rightarrow$  best solution found

create a 2D grid of cells  
for each cell  $i$  in the grid:  
    assign a random value  $x_i$  within  $[lb, ub]$   
for each cell  $i$ :  
     $fitness[i] = f(x_i)$   
    best\_solution = cell with minimum (or max) fitness  
    for iter = 1 to max\_iterations do:  
        for each cell  $i$  in the grid (in parallel):  
            neighbors = cells in neighbourhood of  $i$   
            best\_neighbor = neighbor with best fitness  
             $x_{i\_new} = (x_i + best\_neighbor.value)/2$   
        Enforce boundaries: ensure  $x_{i\_new} \in [lb, ub]$   
        Update all cells simultaneously:  
             $x_i = x_{i\_new}$   
    Recalculate fitness for all cells:  
         $fitness[i] = f(x_i)$   
    Update global best\_solution  
return best\_solution

for  
 b/w 1/25

**Code:**

```
import numpy as np

# Objective function to minimize
def objective_function(x):
    return x**2 - 4*x + 4      # simple convex function, min at x = 2

# Parallel Cellular Algorithm
def parallel_cellular_algorithm(obj_func, grid_size=(10, 10), lb=-10, ub=10,
max_iter=100):
    rows, cols = grid_size
    num_cells = rows * cols

    # Initialize cells randomly in search space
    cells = np.random.uniform(lb, ub, (rows, cols))

    # Evaluate initial fitness
    fitness = obj_func(cells)

    # Define neighborhood (3x3 Moore neighborhood)
    def get_neighbors(r, c):
        neighbors = []
        for i in range(r-1, r+2):
            for j in range(c-1, c+2):
                if (0 <= i < rows) and (0 <= j < cols) and not (i == r and j == c):
                    neighbors.append((i, j))
        return neighbors

    # Track best solution
    best_value = np.min(fitness)
    best_position = np.unravel_index(np.argmin(fitness), fitness.shape)

    # Main iteration loop
    for t in range(max_iter):
        new_cells = np.copy(cells)

        for r in range(rows):
            for c in range(cols):
                neighbors = get_neighbors(r, c)
                # Find the best neighbor
                neighbor_values = np.array([cells[i, j] for i, j in neighbors])
                neighbor_fitness = np.array([obj_func(cells[i, j]) for i, j in neighbors])

                best_neighbor_value = neighbor_values[np.argmin(neighbor_fitness)]

                # Update rule: average of current cell and best neighbor
                new_cells[r, c] = (cells[r, c] + best_neighbor_value) / 2

        # Apply boundaries
        new_cells = np.clip(new_cells, lb, ub)

        # Update cells and fitness
        cells = new_cells
        fitness = obj_func(cells)

    # Track global best
    current_best = np.min(fitness)
```

```

        if current_best < best_value:
            best_value = current_best
            best_position = np.unravel_index(np.argmin(fitness), fitness.shape)

        # Optional: print progress
        if (t + 1) % 10 == 0 or t == 0:
            print(f"Iteration {t+1}/{max_iter} → Best fitness: {best_value:.6f}")

    print("\nBest solution found:")
    print(f"x = {cells[best_position]:.6f}, fitness = {best_value:.6f}")
    return cells[best_position], best_value

# Example run
best_x, best_fit = parallel_cellular_algorithm(
    objective_function,
    grid_size=(10, 10),
    lb=-10,
    ub=10,
    max_iter=50
)

```

Output:

---

```

Iteration 1/50 → Best fitness: 0.000303
Iteration 10/50 → Best fitness: 0.000000
Iteration 20/50 → Best fitness: 0.000000
Iteration 30/50 → Best fitness: 0.000000
Iteration 40/50 → Best fitness: 0.000000
Iteration 50/50 → Best fitness: 0.000000

```

```

Best solution found:
x = 2.000000, fitness = 0.000000

```