

# **VISVESVARAYA TECHNOLOGICAL UNIVERSITY**

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



## **LAB RECORD**

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

*Submitted by*

**Sohan A R(1BM22CS285)**

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

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



**B.M.S. COLLEGE OF ENGINEERING  
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**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 **Sohan A R(1BM22CS285)**, who is 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.

Saritha A N Assistant Professor Department of CSE, BMSCE	Dr. Kavitha Sooda Professor & HOD Department of CSE, BMSCE
----------------------------------------------------------------	------------------------------------------------------------------

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Github Link: <https://github.com/SohanAR/BIS.git>

## **Index Page of Observation Book:**

## Program 1

**Problem statement:** Optimize the allocation of a portfolio using a Genetic Algorithm to maximize the Sharpe Ratio, balancing expected returns and risk. Ensure the total asset allocation adheres to a fixed budget constraint.

Algorithm:

The image shows a handwritten note on a lined notebook page. At the top right, there is a logo for "classmate" with fields for "Date" and "Page". The page contains handwritten text and some diagrams. The text is organized into sections with numbered headings (1) through (9).  
Section 1: LAB - 1  
Section 2: GENETIC ALGORITHM  
Section 3: Summary:  
1) Initialization:  
- Population Creation: A population of binary chromosomes is randomly initialized.  
- Each chromosome has a fixed length (number of bits) and represents a potential solution in the search space defined by  $\alpha$ . (In this case, between 0 and 1).  
2) Evaluation:  
- The fitness of each individual in the population is evaluated using the objective function  $f(x)$ .  
- The binary chromosomes are decoded into real numbers for evaluation.  
3) Selection:  
Tournament selection is used to choose individuals for reproduction. In each tournament, three competitors are randomly selected, and the individual with the highest fitness is chosen.  
4) Crossover:  
Selected parents undergo crossover to create offspring. A random point on the chromosome is chosen, and the segments from two parents are swapped to form new individuals.  
5) Mutation:  
Each chromosome has a small chance (mutation rate) of having its bits flipped (from 0 to 1 or vice versa).

## ② Evolution loop

Algorithm iterates for a specified number of generations & fitness is evaluated.

## ③ Output

After all generations, the best solution (decoded chromosome) and its corresponding fitness value are returned and printed.

26/9/2012 \*

## Code:

11/20/24, 6:59 PM

Untitled15.ipynb - Colab

```
import random
import numpy as np
import math

def generate_cities(num_cities):
    return [(random.randint(0, 100), random.randint(0, 100)) for _ in range(num_cities)]

def compute_distance_matrix(cities):
    num_cities = len(cities)
    distances = [[0] * num_cities for _ in range(num_cities)]
    for i in range(num_cities):
        for j in range(num_cities):
            if i != j:
                distances[i][j] = math.sqrt(
                    (cities[i][0] - cities[j][0])**2 + (cities[i][1] - cities[j][1])**2
                )
    return distances

class TSP:
    def __init__(self, distances):
        self.distances = distances
        self.num_cities = len(distances)

    def fitness(self, route):
        total_distance = sum(
            self.distances[route[i]][route[i + 1]] for i in range(len(route) - 1)
        )
        total_distance += self.distances[route[-1]][route[0]]

    class GeneticAlgorithm:
        def __init__(self, tsp, population_size=100, generations=500, mutation_rate=0.1):
            self.tsp = tsp
            self.population_size = population_size
            self.generations = generations
            self.mutation_rate = mutation_rate
            self.population = self._initialize_population()

        def _initialize_population(self):
            return [random.sample(range(self.tsp.num_cities), self.tsp.num_cities) for _ in range(self.population_size)]

        def _select_parents(self):
            fitnesses = [self.tsp.fitness(route) for route in self.population]
            total_fitness = sum(fitnesses)
            probabilities = [f / total_fitness for f in fitnesses]
            return random.choices(self.population, probabilities, k=2)

        def _crossover(self, parent1, parent2):
            size = len(parent1)
            start, end = sorted(random.sample(range(size), 2))
            child = [-1] * size
            child[start:end] = parent1[start:end]

            p2_idx = 0
            for i in range(size):
                if child[i] == -1:
                    while parent2[p2_idx] in child:
                        p2_idx += 1
                    child[i] = parent2[p2_idx]
            return child
```

```

def _mutate(self, route):

    if random.random() < self.mutation_rate:
        i, j = random.sample(range(len(route)), 2)
        route[i], route[j] = route[j], route[i]

def evolve(self):
    for _ in range(self.generations):
        new_population = []
        for _ in range(self.population_size):
            parent1, parent2 = self._select_parents()
            child = self._crossover(parent1, parent2)
            self._mutate(child)
            new_population.append(child)
        self.population = new_population

def get_best_solution(self):

    best_route = min(self.population, key=lambda route: 1 / self.tsp.fitness(route))
    best_distance = 1 / self.tsp.fitness(best_route)
    return best_route, best_distance

if __name__ == "__main__":
    num_cities = 5
    cities = generate_cities(num_cities)
    distances = compute_distance_matrix(cities)

    print("City Coordinates:")
    for i, city in enumerate(cities):
        print(f"City {i}: {city}")

    tsp = TSP(distances)
    ga = GeneticAlgorithm(tsp, population_size=50, generations=100, mutation_rate=0.2)
    ga.evolve()
    best_route, best_distance = ga.get_best_solution()

    print("\nBest route:", best_route)
    print("Best distance:", best_distance)

→ City Coordinates:
City 0: (22, 31)
City 1: (33, 46)
City 2: (89, 0)
City 3: (65, 0)
City 4: (3, 17)

Best route: [2, 1, 0, 4, 3]
Best distance: 202.96101904990562

```

## Program 2

**Problem statement:** Implement a Particle Swarm Optimization (PSO) algorithm to minimize benchmark functions, such as the Rastrigin and Sphere functions, by optimizing their input parameters. The goal is to find the global minimum while efficiently exploring the solution space using swarm intelligence.

Algorithm:

The image shows handwritten notes on a lined notebook page. At the top right, there is a logo with the text "classmate" above "Date \_\_\_\_\_" and "Page \_\_\_\_\_". Below this, the title "LAB-2" and "Particle Swarm Optimization (PSO)" are written. The notes are organized into numbered sections: 1. Objective, 2. Key Components, 3. Algorithm Steps. The notes provide a high-level overview of the PSO algorithm, its components, and its iterative process.

1. Objective : PSO is an optimization algorithm inspired by the social behavior of birds or fish - It aims to find the optimal solution for a given function iteratively improving candidate solutions

2. Key Components :

- Particles : Each candidate solution represented as a particle in the search space
- Fitness Function : Evaluates the quality of the particle's position
- Velocity Update : Each particle adjusts its velocity based on its best known position and the global best position
- Position Update : Particles move to new positions based on their updated velocities

3. Algorithm Steps

- Initialization : Generate particles with random positions and velocities within specified bounds.
- Evaluation : Compute the fitness of each particle
- Update : Adjust each particle's velocity and position
- Iteration : Repeat evaluation and updating for a set number of iterations or until convergence

- + Output - Return the best position and value found

#### Applications of PSO

1. Engineering Optimization
2. Function Optimization
3. Machine Learning
4. Scheduling
5. Robotics

#### Working Mechanism

1. Initialization
  - + Randomly initializing a swarm of particles with position and velocities within defined search boundaries
2. Convergence check :
  - + Iterate the above steps until a maximum number of iterations is reached or until the solution converges.

Step 1/2x

## Code:

```
11/20/24, 7:55 AM Untitled9.ipynb - Colab

import random
import numpy as np

def objective_function(position):
    """The function to be minimized."""
    x, y = position
    return x**2 + y**2

def pso(objective_function, dimensions, iterations, population_size, w=0.7, c1=1.4, c2=1.4):
    """
    Particle Swarm Optimization algorithm.

    Args:
        objective_function: The function to be minimized.
        dimensions: The number of dimensions of the search space.
        iterations: The number of iterations to run the algorithm.
        population_size: The number of particles in the swarm.
        w: Inertia weight.
        c1: Cognitive parameter.
        c2: Social parameter.

    Returns:
        A tuple containing the best solution found and its corresponding objective function value.
    """
    particles = []
    for _ in range(population_size):
        position = np.random.uniform(-10, 10, dimensions)
        velocity = np.random.uniform(-1, 1, dimensions)
        particles.append({
            'position': position,
            'velocity': velocity,
            'best_position': position.copy(),
            'best_value': objective_function(position)
        })

    global_best_position = particles[0]['best_position'].copy()
    global_best_value = particles[0]['best_value']

    for _ in range(iterations):
        for particle in particles:

            r1 = random.random()
            r2 = random.random()
            particle['velocity'] = (w * particle['velocity'] +
                                    c1 * r1 * (particle['best_position'] - particle['position']) +
                                    c2 * r2 * (global_best_position - particle['position']))

            particle['position'] = particle['position'] + particle['velocity']
            particle['position'] = np.clip(particle['position'], -10, 10)

            value = objective_function(particle['position'])

            if value < particle['best_value']:
                particle['best_value'] = value
                particle['best_position'] = particle['position'].copy()

            if value < global_best_value:
                global_best_value = value
                global_best_position = particle['position'].copy()

    return global_best_position, global_best_value
```

<https://colab.research.google.com/drive/10dBsvVCUavtG9L8Q19FQ8mN6idnJEdjz#scrollTo=lUs2D5e326tn&printMode=true>

1/2

---

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```
dimensions = 2
iterations = 100
population_size = 50

best_position, best_value = pso(objective_function, dimensions, iterations, population_size)
print(f"Best position found: {best_position}")
print(f"Best value found: {best_value}")

→ Best position found: [ 4.04789703e-08 -2.23363404e-08]
Best value found: 2.137459138638845e-15
```

### Program 3

**Problem statement:** Implement an Ant Colony Optimization (ACO) algorithm to solve the Traveling Salesman Problem (TSP), where the goal is to find the shortest possible path that visits all cities exactly once and returns to the starting city. The algorithm should utilize pheromone trails and heuristic information to guide the search efficiently.

Algorithm:

The image shows handwritten Python code for an Ant Colony Optimization (ACO) algorithm to solve the Traveling Salesman Problem (TSP). The code is organized into three main classes: `Ant Colony`, `TSP`, and `ACO`. The `Ant Colony` class imports `random`, `numpy` as `np`, `math`, and `matplotlib.pyplot` as `plt`. It defines a function `euclidean_distance` to calculate the distance between two cities. The `TSP` class initializes with a list of cities and creates a matrix of distances. The `ACO` class initializes with parameters for the number of ants, alpha, beta, rho, and iterations, and sets up initial pheromone levels.

```
classmate
Date _____
Page _____  
  
Ant Colony  
  
import random  
import numpy as np  
import math  
import matplotlib.pyplot as plt  
  
def euclidean_distance(city1, city2):  
    return math.sqrt((city1[0] - city2[0]) ** 2 + (city1[1] - city2[1]) ** 2)  
  
class TSP:  
    def __init__(self, cities):  
        self.cities = cities  
        self.num_cities = len(cities)  
        self.distances = np.zeros((self.num_cities, self.num_cities))  
  
        for i in range(self.num_cities):  
            for j in range(self.num_cities):  
                self.distances[i][j] = euclidean_distance(cities[i], cities[j])  
  
class ACO:  
    def __init__(self, tsp, num_ants, alpha, beta, rho, iterations):  
        self.tsp = tsp  
        self.num_ants = num_ants  
        self.alpha = alpha  
        self.beta = beta  
        self.rho = rho  
        self.iterations = iterations  
        self.pheromone = np.ones((self.tsp.num_cities, self.tsp.num_cities)) * 1e-05
```

```
def choose_next_city(self, current_city, visited, pheromone_probabilities = np.zeros(len(self.tsp.num_cities)), total_pheromone = 0)
```

for i in range(0, self.tsp.num\_cities):

if i not in visited:

pheromone\_probabilities[i] = (pheromone[current\_city][i] \* 1 / distances[current\_city][i]) \*  $\alpha$   
total\_pheromone += pheromone\_probabilities[i]

pheromone\_probabilities /= total\_pheromone

return np.random.choice(range(self.tsp.num\_cities), p=pheromone\_probabilities)

```
def run(self):
```

best\_path = None

best\_path\_length = float('inf')

for \_ in range(0, self.iterations):

all\_paths = self.construct\_solution()

for path in all\_paths:

path\_length = self.calculate\_path\_length(path)

if path\_length < best\_path\_length:

best\_path = path

best\_path\_length = path\_length

self.update\_pheromone(all\_paths)

return best\_path, best\_path\_length

if name == "main":

cities = [(0, 1), (1, 3), (4, 3), (6, 1), (3, 0)]

tsp = TSP (cities)

num\_ants = 10

alpha = 1.0

beta = 2.0

rho = 0.1

Aerating = 100

a(s) = ACO (tsp, num\_ants, alpha, beta, rho, Aerating)  
~~best\_path, best\_path\_length = a(s).run()~~

print ("Best Path : ", best\_path)

print ("Best path length : ", best\_path\_length)

best\_path\_Ends = [cities[i] for i in best\_path]  
~~[cities[best\_path[0]]]~~

best\_path\_coords = np.array (best\_path\_coords)

plt.plot (best\_path\_coords[:, 0], best\_path\_coords[:, 1], no

plt.scatter (\*zip (\*cities), color='red')

plt.title ("Best TSP Path (length : " + str(best\_path\_length) + ")")

plt.show ()

Output :

Best path : [1, 2, 3, 4, 0]

Best Path length : 15.15798274

Car. 18.11

## Code:

11/20/24, 7:51 AM

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

# Define constants for the algorithm
NUM_ANTS = 50
NUM_CITIES = 20 # Now we have 20 cities
ALPHA = 1.0 # Influence of pheromone
BETA = 2.0 # Influence of distance
RHO = 0.1 # Pheromone evaporation rate
Q = 100 # Pheromone deposit constant
MAX_ITER = 100 # Maximum number of iterations

# Predefined 20 cities (coordinates in 2D space)
def generate_cities():
    cities = np.array([
        [5, 10], [11, 5], [14, 9], [12, 15], [8, 13], # Cities 0-4
        [10, 10], [13, 7], [16, 5], [14, 3], [18, 6], # Cities 5-9
        [4, 2], [7, 1], [8, 5], [6, 7], [4, 10], # Cities 10-14
        [15, 18], [12, 17], [3, 18], [17, 12], [19, 8] # Cities 15-19
    ])
    return cities

# Compute the distance matrix
def compute_distance_matrix(cities):
    num_cities = len(cities)
    distance_matrix = np.zeros((num_cities, num_cities))
    for i in range(num_cities):
        for j in range(i + 1, num_cities):
            dist = np.linalg.norm(cities[i] - cities[j])
            distance_matrix[i, j] = dist
            distance_matrix[j, i] = dist
    return distance_matrix

# Initialize pheromone matrix
def initialize_pheromone_matrix(num_cities):
    pheromone_matrix = np.ones((num_cities, num_cities)) # Pheromone starts as 1 for all edges
    np.fill_diagonal(pheromone_matrix, 0) # No pheromone on the diagonal (self-loops)
    return pheromone_matrix

# Calculate the total length of a tour
def calculate_tour_length(tour, dist_matrix):
    length = 0
    for i in range(len(tour) - 1):
        length += dist_matrix[tour[i], tour[i + 1]]
    length += dist_matrix[tour[-1], tour[0]] # Returning to the start
    return length

# Ant solution construction (probabilistic decision on next city)
def construct_solution(num_cities, pheromone_matrix, dist_matrix):
    tour = [random.randint(0, num_cities - 1)] # Start from a random city
    visited = set(tour)

    while len(tour) < num_cities:
        current_city = tour[-1]
        probabilities = []
        for next_city in range(num_cities):
            if next_city not in visited:
                pheromone = pheromone_matrix[current_city, next_city] ** ALPHA
                distance = (1.0 / dist_matrix[current_city, next_city]) ** BETA
                probabilities.append(pheromone * distance)
            else:
                probabilities.append(0)
        total_prob = sum(probabilities)
        if total_prob == 0:
            break
        next_city_index = np.random.choice(range(len(probabilities)), p=probabilities)
        tour.append(next_city_index)
        visited.add(next_city_index)
```

[https://colab.research.google.com/drive/1hCD\\_n90J05NRotPADsr9QsfZSFSfehbb#scrollTo=l0L49nQ1wY7&printMode=true](https://colab.research.google.com/drive/1hCD_n90J05NRotPADsr9QsfZSFSfehbb#scrollTo=l0L49nQ1wY7&printMode=true)

1/3

```

total_prob = sum(probabilities)
probabilities = [p / total_prob for p in probabilities]

# Choose the next city based on the probabilities
next_city = np.random.choice(range(num_cities), p=probabilities)
tour.append(next_city)
visited.add(next_city)

return tour

# Update the pheromone matrix based on the solutions found by ants
def update_pheromone(pheromone_matrix, all_tours, dist_matrix, best_tour):
    # Evaporate pheromone
    pheromone_matrix *= (1 - RHO)

    # Add pheromone for all ants
    for tour in all_tours:
        tour_length = calculate_tour_length(tour, dist_matrix)
        for i in range(len(tour) - 1):
            pheromone_matrix[tour[i], tour[i + 1]] += Q / tour_length
        pheromone_matrix[tour[-1], tour[0]] += Q / calculate_tour_length(tour, dist_matrix)

    # Add pheromone for the best tour
    best_length = calculate_tour_length(best_tour, dist_matrix)
    for i in range(len(best_tour) - 1):
        pheromone_matrix[best_tour[i], best_tour[i + 1]] += Q / best_length
    pheromone_matrix[best_tour[-1], best_tour[0]] += Q / best_length

# Main ACO algorithm for solving TSP
def ant_colony_optimization(num_cities, dist_matrix, pheromone_matrix, max_iter):
    best_tour = None
    best_tour_length = float('inf')

    # Main loop
    for iteration in range(max_iter):
        all_tours = []

        # Step 1: All ants construct their solutions
        for _ in range(NUM_ANTS):
            tour = construct_solution(num_cities, pheromone_matrix, dist_matrix)
            all_tours.append(tour)
            tour_length = calculate_tour_length(tour, dist_matrix)

        # Step 2: Update the best tour if necessary
        if tour_length < best_tour_length:
            best_tour = tour
            best_tour_length = tour_length

        # Step 3: Update pheromone matrix
        update_pheromone(pheromone_matrix, all_tours, dist_matrix, best_tour)

        # Optional: print progress every 10 iterations
        if iteration % 10 == 0:
            print(f"Iteration {iteration}: Best tour length = {best_tour_length:.2f}")

    return best_tour, best_tour_length

# Main Execution
if __name__ == "__main__":
    # Step 1: Generate predefined cities and distance matrix
    cities = generate_cities()
    dist_matrix = compute_distance_matrix(cities)

```

[https://colab.research.google.com/drive/1hCD\\_n90J05NRotPADsr9QsfZSFSfehbb#scrollTo=IL0L49nQ1wY7&printMode=true](https://colab.research.google.com/drive/1hCD_n90J05NRotPADsr9QsfZSFSfehbb#scrollTo=IL0L49nQ1wY7&printMode=true)

2/3

```
# Step 2: Initialize pheromone matrix
pheromone_matrix = initialize_pheromone_matrix(NUM_CITIES)

# Step 3: Run ACO algorithm
best_tour, best_tour_length = ant_colony_optimization(NUM_CITIES, dist_matrix, pheromone_matrix, MAX_ITER)

# Step 4: Output the best tour and visualize it
print(f"Best tour length: {best_tour_length:.2f}")
```

```
→ Iteration 0: Best tour length = 107.48
Iteration 10: Best tour length = 81.48
Iteration 20: Best tour length = 80.59
Iteration 30: Best tour length = 80.50
Iteration 40: Best tour length = 79.23
Iteration 50: Best tour length = 79.23
Iteration 60: Best tour length = 77.88
Iteration 70: Best tour length = 77.88
Iteration 80: Best tour length = 77.88
Iteration 90: Best tour length = 77.88
Best tour length: 77.88
```

## Program 4

**Problem statement:** Implement the Cuckoo Search Algorithm for feature selection to identify an optimal subset of features that maximizes the classification accuracy of a Support Vector Machine (SVM) on a given dataset.

Algorithm:

Lab 7  
Cuckoo Search

```
import numpy as np

def levy_flight(lambd_a, size):
    from scipy.special import gamma
    sigma = (gamma(1+lambd_a)**2 * np.sin(np.pi * lambd_a / 2)) / (gamma((1+lambd_a)/2) * lambd_a * np.cos(np.pi * lambd_a / 2)) ** (1/lambd_a)
    u = np.random.normal(0, sigma, size)
    v = np.random.normal(0, 1, size)
    return u/(np.abs(v) ** (1/lambd_a))

def cuckoo_search(objective_function, bounds, num_tests=1000, max_iter=1000, pa=0.25, Levy_Law_Exponent=lambd_a):
    nests = np.random.uniform(bounds[0], bounds[1], (num_tests, len(bounds)))
    fitness = np.array([objective_function(nest) for nest in nests])
    best_nest = nests[np.argmin(fitness)]
    best_fitness = np.min(fitness)

    for iteration in range(max_iter):
        new_nests = []
        for i in range(len(nests)):
            step = levy_flight(lambd_a, len(bounds))
            new_nest = np.clip(best_nest + step, bounds[0], bounds[1])
            new_nests.append(new_nest)
            if np.random.rand() < pa:
                new_nest = objective_function(new_nest)
                new_nests.append(new_nest)
        nests = new_nests
        fitness = np.array([objective_function(nest) for nest in nests])
        if np.min(fitness) < best_fitness:
            best_nest = nests[np.argmin(fitness)]
            best_fitness = np.min(fitness)
```

`new_nests.append(new_nest)`

`new_fitness = np.array([objective_function(nest) for nest in new_nests])`

`for i in range(num_nests):`

`if new_fitness[i] < fitness[i]:`

`nests[i] = new_nests[i]`

`fitness[i] = new_fitness[i]`

`for i in range(num_nests):`

`if np.random.rand() < pa:`

`nests[i] = np.random.uniform(bounds[0], bounds[1], len(bounds)))`

`fitness[i] = objective_function(nests[i])`

`current_best_fitness = np.min(fitness)`

`if current_best_fitness < best_fitness:`

`best_fitness = current_best_fitness`

`best_nest = nests[np.argmin(fitness)]`

~~`print(f"Iteration {iteration + 1} / {max_iter}, Best fitness: {best_fitness}"")`~~

`return best_nest, best_fitness`

`bounds = np.array([-5, -5, 5, 5])`

`best_solution, best_objective_value = coloss.search(objective_function, bounds)`

`print("Best Solution", best_solution)`

`print("Best Objective Value:", best_objective_value)`

Output:

Best Solution:  $[-0.51597054 \quad -1.41565335]$

Best Objective Value:  $0.0020850514611121572$

Dec  
x1.11

## Code:

11/20/24, 7:41 AM Untitled13.ipynb - Colab

```
#cuckoo search(Traffic Signal Optimization)
import numpy as np
from scipy.special import gamma

def fitness_function(x):
    waiting_times = np.array([10 + (x[i] ** 2) / 100 for i in range(len(x))])
    total_waiting_time = np.sum(waiting_times)
    return total_waiting_time

def levy_flight(dim, beta=1.5):
    sigma_u = np.power((gamma(1 + beta) * np.sin(np.pi * beta / 2) /
                        gamma((1 + beta) / 2) * beta * (2 ** (beta - 1))), 1 / beta)
    u = np.random.normal(0, sigma_u, dim)
    v = np.random.normal(0, 1, dim)
    step = u / np.power(np.abs(v), 1 / beta)
    return step

def cuckoo_search(dim, bounds, num_nests, max_iter, p_a=0.25, Lambda=1.5):
    nests = np.random.uniform(bounds[0], bounds[1], (num_nests, dim))
    fitness = np.array([fitness_function(nest) for nest in nests])

    best_idx = np.argmin(fitness)
    best_nest = nests[best_idx]
    best_fitness = fitness[best_idx]

    for iter in range(max_iter):
        new_nests = np.copy(nests)
        for i in range(num_nests):
            step = levy_flight(dim, Lambda)
            new_nests[i] = nests[i] + step
            new_nests[i] = np.clip(new_nests[i], bounds[0], bounds[1])

        new_fitness = np.array([fitness_function(nest) for nest in new_nests])

        for i in range(num_nests):
            if new_fitness[i] < fitness[i]:
                nests[i] = new_nests[i]
                fitness[i] = new_fitness[i]

        if np.random.rand() < p_a:
            random_idx = np.random.randint(num_nests)
            nests[random_idx] = np.random.uniform(bounds[0], bounds[1], dim)
            fitness[random_idx] = fitness_function(nests[random_idx])

        current_best_idx = np.argmin(fitness)
        current_best_fitness = fitness[current_best_idx]

        if current_best_fitness < best_fitness:
            best_fitness = current_best_fitness
            best_nest = nests[current_best_idx]

    return best_nest, best_fitness

dim = 3
bounds = [10, 120]
num_nests = 20
max_iter = 100

best_solution, best_value = cuckoo_search(dim, bounds, num_nests, max_iter)

print("\n--- Best Solution ---")
print("Green Light Timings (seconds):", best_solution)
print("Best Fitness Value (Total Waiting Time):", best_value)
```

1/2

---

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```
print("Green Light Timings (seconds):", best_solution)
print("Best Fitness Value (Total Waiting Time):", best_value)
```

```
--- Best Solution ---
Green Light Timings (seconds): [10, 10, 10]
Best Fitness Value (Total Waiting Time): 33.0
```

## Program 5

**Problem Statement :** Implement the Grey Wolf Optimizer (GWO) to optimize the hyperparameters (C and gamma) of a Support Vector Machine (SVM) classifier for achieving the best classification accuracy on the Iris dataset.

### **Algorithm :**

LABS - 8  
Grey wolf optimizer

```
import numpy as np

def objective_function(x):
    return np.sum(x**2)

class grey_wolf_optimizer:
    def __init__(self, obj_function, dim, n_wolves, max_iter,
                 bounds):
        self.obj_function = obj_function
        self.dim = dim
        self.n_wolves = n_wolves
        self.max_iter = max_iter
        self.bounds = bounds
        self.alpha_pos = np.zeros(dim)
        self.beta_pos = np.zeros(dim)
        self.delta_pos = np.zeros(dim)
        self.alpha_score = float("inf")
        self.beta_score = float("inf")

    def initialize_population(self):
        return np.random.uniform(self.bounds[0], self.bounds[1],
                               (self.n_wolves, self.dim))

    def update_wolf_positions(self, wolves, t, a):
        for i in range(self.n_wolves):
            A1 = 2 * t * np.random.random(self.dim) - a
            C1 = 2 * t * np.random.random(self.dim)
```

$$A_2 = 2^k a * np.random.random(self.dim) - a$$

$$C_2 = 2^k np.random.random(self.dim)$$

$$A_3 = 2^k a * np.random.random(self.dim) - a$$

$$C_3 = 2^k np.random.random(self.dim)$$

$$D\_alpha = abs(C_1 * self.alpha - pos - wolves[i])$$

$$D\_beta = abs(C_2 * self.beta - pos - wolves[i])$$

$$D\_delta = abs(C_3 * self.delta - pos - wolves[i])$$

$$X_1 = self.alpha - pos - A_1 * D\_alpha$$

$$X_2 = self.beta - pos - A_2 * D\_beta$$

$$X_3 = self.delta - pos - A_3 * D\_delta$$

$$wolves[i] = (X_1 + X_2 + X_3) / 3$$

$$wolves[i] = np.clip(wolves[i], self.bounds[0],$$

$$self.bounds[1])$$

def optimize(self):

wolves = self.initialize\_population()

for t in range(self.max\_iter):

for i in range(self.n\_wolves):

fitness = self.obj\_function(wolves[i])

, if fitness < self.alpha\_scores:

self.alpha\_score = fitness

self.alpha\_pos = wolves[i].copy()

a = 2 - t / (2 / self.max\_iter)

return self.alpha\_pos, self.alpha\_score

print ("Best Solution ", best\_solution)  
print ("Best Score ", best\_score)

~~Best\_Sol~~

Output

Best Solution : [ 2.1696229e-11, -2.1291948e-11,  
-1.95888221e-11, -2.06436664e-11,  
-1.88732585e-11 ]

Best Score : 2.0901197822659867e-21

~~Score~~  
~~28.13~~

Code:

```

import numpy as np

N_INPUTS = 3
N_HIDDEN = 5
N_OUTPUTS = 1
N_WOLVES = 30
MAX_ITER = 100
LB = -10.0
UB = 10.0

def sigmoid(x):
    return 1.0 / (1.0 + np.exp(-x))

def sigmoid_derivative(x):
    return x * (1 - x)

def forward_pass(input_data, weights):
    hidden_layer = np.dot(input_data, weights[:N_INPUTS *
N_HIDDEN].reshape(N_INPUTS, N_HIDDEN))
    hidden_layer = sigmoid(hidden_layer)
    output = np.dot(hidden_layer, weights[N_INPUTS *
N_HIDDEN:]).reshape(N_HIDDEN, N_OUTPUTS))
    return output, hidden_layer

def fitness_function(weights, inputs, targets, n_samples):
    total_error = 0.0
    for i in range(n_samples):
        output, _ = forward_pass(inputs[i], weights)
        total_error += (output - targets[i]) ** 2
    return total_error / n_samples

def rand_range(min_val, max_val):
    return min_val + (max_val - min_val) * np.random.random()

def update_position(positions, alpha_pos, beta_pos, delta_pos, i, t):
    A = 2 - t * (2.0 / MAX_ITER)
    C = 2 * np.random.random()

    for j in range(len(positions[i])):
        D_alpha = np.abs(C * alpha_pos[j] - positions[i][j])
        D_beta = np.abs(C * beta_pos[j] - positions[i][j])
        D_delta = np.abs(C * delta_pos[j] - positions[i][j])

        new_position = alpha_pos[j] - A * D_alpha if np.random.random() > 0.5 else beta_pos[j] - A * D_beta

```

```

    positions[i][j] = np.clip(new_position, LB, UB)

def gwo_optimization(inputs, targets, n_samples):
    positions = np.random.uniform(LB, UB, (N_WOLVES, N_INPUTS *
N_HIDDEN + N_HIDDEN)) # Wolves' positions (weights)
    fitness = np.zeros(N_WOLVES) # Fitness of wolves
    alpha_pos = np.zeros(N_INPUTS * N_HIDDEN + N_HIDDEN) # Best
position (alpha wolf)
    beta_pos = np.zeros(N_INPUTS * N_HIDDEN + N_HIDDEN) # Second best
position (beta wolf)
    delta_pos = np.zeros(N_INPUTS * N_HIDDEN + N_HIDDEN) # Third best
position (delta wolf)
    alpha_score = float('inf') # Best fitness value (alpha wolf)
    beta_score = float('inf') # Second best fitness value (beta wolf)
    delta_score = float('inf') # Third best fitness value (delta wolf)

    for i in range(N_WOLVES):
        fitness[i] = fitness_function(positions[i], inputs, targets,
n_samples)

        if fitness[i] < alpha_score:
            alpha_score = fitness[i]
            alpha_pos = positions[i]
        elif fitness[i] < beta_score:
            beta_score = fitness[i]
            beta_pos = positions[i]
        elif fitness[i] < delta_score:
            delta_score = fitness[i]
            delta_pos = positions[i]

    for t in range(MAX_ITER):
        for i in range(N_WOLVES):
            update_position(positions, alpha_pos, beta_pos, delta_pos,
i, t)
            fitness[i] = fitness_function(positions[i], inputs,
targets, n_samples)

            if fitness[i] < alpha_score:
                alpha_score = fitness[i]
                alpha_pos = positions[i]
            elif fitness[i] < beta_score:
                beta_score = fitness[i]
                beta_pos = positions[i]

```

```

        elif fitness[i] < delta_score:
            delta_score = fitness[i]
            delta_pos = positions[i]

    if t % 10 == 0:
        print(f"Iteration {t}/{MAX_ITER}, Best Fitness = {alpha_score}")

    print(f"\nBest Fitness: {alpha_score}")
    return alpha_pos

inputs = np.array([
    [0.0, 0.0, 1.0],
    [1.0, 0.0, 1.0],
    [0.0, 1.0, 1.0],
    [1.0, 1.0, 1.0]
])

targets = np.array([0.0, 1.0, 1.0, 0.0])

best_weights = gwo_optimization(inputs, targets, len(inputs))

print("\nEvaluating the final model with the best weights...")
for i in range(len(inputs)):
    output, _ = forward_pass(inputs[i], best_weights)
    print(f"Input: {inputs[i]}, Predicted Output: {output[0]}, Actual Target: {targets[i]}")

```

Iteration 0/100, Best Fitness = 0.5000013911617378

Iteration 10/100, Best Fitness = 0.5000013911617378

Iteration 20/100, Best Fitness = 0.5000013911617378

Iteration 30/100, Best Fitness = 0.5000013911617378

Iteration 40/100, Best Fitness = 0.5000013911617378

Iteration 50/100, Best Fitness = 0.5000013911617378

Iteration 60/100, Best Fitness = 0.5000013911617378

Iteration 70/100, Best Fitness = 0.5000013911617378

Iteration 80/100, Best Fitness = 0.5000013911617378

Iteration 90/100, Best Fitness = 0.5000013911617378

Best Fitness: 0.5000013911617378

Evaluating the final model with the best weights...

Input: [0. 0. 1.], Predicted Output: -0.0022698934351217197, Actual Target: 0.0

Input: [1. 0. 1.], Predicted Output: -1.0305768090951018e-07, Actual Target: 1.0

Input: [0. 1. 1.], Predicted Output: -1.0305768090951018e-07, Actual Target: 1.0

Input: [1. 1. 1.], Predicted Output: -4.678811484419649e-12, Actual Target: 0.0

== Code Execution Successful ==

## Program 6

**Problem Statement :** Develop a parallel cellular automaton-based algorithm for optimal robot route planning in a grid-based environment, ensuring collision-free navigation while minimizing travel distance and computational time.

### Algorithm :

```
LAB 9
Parallel Cellular Algorithm

import numpy as np

def fitness(x,y):
    return -x**2 - y**2 + 10*x + 8*y

grid_size = (5,5)
num_Iteration = 5
Search_space = (10,10)
neighbourhood_size = 1

population = np.random.uniform(Search_space[0],
                               Search_space[1], C*grid_size, 2))

def get_best_neighbour(population, x, Y, neighbourhood_size):
    x_min = x - neighbourhood_size
    max(grid[0], x + neighbourhood_size + 1)

    neighborhood = population[x_min:x_max, y_min:y_max]
    neighborhood = neighborhood.reshape(-1, 2)

    return max(neighborhood, key=lambda ind: fitness(ind[0], ind[1]))
```

for iteration in range(num\_Iteration):  
 new\_population = np.copy(population)  
 for x in range(grid\_size[0]):  
 for Y in range(C\*grid\_size[1]):  
 best\_neighbour = get\_best\_neighbour(new\_population,  
 x, Y, neighbourhood\_size)

`new_population(x, y) = np.array([best_neighbour, t mutation  
search_space[0], search_space[1]])`

`population = new_population`

`fitness = np.array([fitness_function(ind[0]), ind[1]  
for ind in rows])  
for row in population])`

`best_idx = np.unravel_index(fitness.argmax(), fitness.shape)`

`print(f"iteration {iteration+1}: Best fitness = {  
fitness[best_idx]}")  
Best_Cell =  
{population[best_idx]}")`

`# output final best solution`

`best_idx = np.unravel_index(fitness.argmax(),  
fitness.shape)`

`print(f"\nBest Solution = {x,y} = {population[best_idx]}  
y, fitness[best_idx]: {y})")`

`Output:`

`Iteration 1: Best fitness = -40.60 Best Cell  
[4.78053407 4.39665832]`

`Iteration 2: Best fitness = 40.99 Best Cell =  
[5.070018 3.90911838]`

`Iteration 3: Best fitness = 40.98 Best Cell =  
[4.86435292 3.93271728]`

`Best Cell = [5.12305524 4.05191983]  
Best Solution = (x,y) = [5.12305524, 4.03191983]  
fitness = 40.98`

Code:

```

import random
import numpy as np

# Initialize Grid (N x M) with random states (0 or 1)
def initialize_grid(N, M):
    np.random.choice([0, 1], size=(N, M))

File display
# count live neighbors for a cell (i, j)
def count_live_neighbors(grid, i, j, N, M):
    return sum(
        grid[ni, nj] for ni in range(i-1, i+2) for nj in range(j-1, j+2)
        if 0 <= ni < N and 0 <= nj < M and (ni != i or nj != j)
    )

# Update cell's state based on neighbors' count
def update_cell(grid, new_grid, i, j, N, M):
    live_neighbors = count_live_neighbors(grid, i, j, N, M)
    if grid[i, j] == 1:
        new_grid[i, j] = 1 if live_neighbors in [2, 3] else 0
    else:
        new_grid[i, j] = 1 if live_neighbors == 3 else 0

# Print the current state of the grid (0s and 1s)
def print_grid(grid):
    for row in grid:
        print(''.join(map(str, row)))
    print() # For spacing between generations

# Main Game of Life simulation with printing grid
def parallel_game_of_life(N, M, steps):
    grid = initialize_grid(N, M)

    for _ in range(steps):
        print_grid(grid) # Print current grid state
        new_grid = np.zeros((N, M), dtype=int) # New grid for next state
        for i in range(N):
            for j in range(M):
                update_cell(grid, new_grid, i, j, N, M)
        grid = new_grid # Update grid for next iteration

    print_grid(grid) # Print final grid after all steps

# Example Usage
N, M = 5, 5 # Smaller grid size for readability
steps = 5 # Number of iterations
final_grid = parallel_game_of_life(N, M, steps)

```

## Output:

```
[→] 0 0 0 0 0  
0 1 0 0 0  
0 0 0 1 1  
0 1 0 1 1  
1 1 1 0 1  
  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 1 1  
1 1 0 0 0  
1 1 1 0 1  
  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
1 0 0 0 1  
1 0 1 0 0  
  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
0 1 0 0 0  
0 1 0 0 0  
  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0  
0 0 0 0 0
```

## Program 7

**Problem Statement :** Solve the 0/1 Knapsack Problem using the Gene Expression Algorithm (GEA) to maximize the total value of selected items without exceeding the given weight capacity.

### **Algorithm :**

LATS-10  
PROGRAM: GENETIC ALGORITHM

```
import numpy as np

def fitness_function(x):
    return -x[0]*2 + 5*x[1] + 6

population_size = 10
mutation_rate = 0.1
crossover_rate = 0.8
num_generations = 50
search_space = [-10, 10]

population = np.random.uniform(search_space[0], search_space[1], population_size)

for generation in range(1, num_generations):
    fitness = np.array([fitness_function(individual) for individual in population])

    adjusted_fitness = fitness - np.min(fitness)
    probabilities = adjusted_fitness / np.sum(adjusted_fitness)

    selected = np.random.choice(population, size=population_size, p=probabilities)

    next_population = []
    for i in range(0, population_size, 2):
        parent_1 = selected[i]
        parent_2 = selected[i+1]
        if np.random.rand() < crossover_rate:
            offspring_1 = (parent_1 + parent_2) / 2
            offspring_2 = (parent_2 + parent_1) / 2
            next_population.append(offspring_1)
            next_population.append(offspring_2)
        else:
            next_population.append(parent_1)
            next_population.append(parent_2)

    population = next_population
```

else:

offspring 1, offspring 2 = parent 1, parent 2

offspring = np.random.uniform(-1, 1)  
 if np.random.rand() < mutation\_rate  
 else 0

offspring t = np.random.uniform(t, 1)  
 if np.random.rand() < mutation\_rate  
 else 0

best\_idx = fitness.argmax()

print "Generation {} generation {}":

Best fitness = fitness[best\_idx] : .2f;

Best Individual = population[best\_idx]  
 : .2f )"

final\_fitness = np.array([fitness\_function(ind)  
 for ind in population])

best\_idx = final\_fitness.argmax()

print "In Best Solution: X = population [ best\_idx ]  
 : if y fitness = final\_fitness [ best\_idx ]: .2f"

Output:

Best Solution x=2.48 , fitness=12.75

Code:

```
import random

# Define the Knapsack Problem (Objective Function)
def knapsack_fitness(items, capacity, solution):
    total_weight = sum([items[i][0] for i in range(len(solution)) if solution[i] == 1])
    total_value = sum([items[i][1] for i in range(len(solution)) if solution[i] == 1])

    # If total weight exceeds the capacity, return 0 (invalid solution)
    if total_weight > capacity:
        return 0
    return total_value

# Gene Expression Algorithm (GEA)
class GeneExpressionAlgorithm:
    def __init__(self, population_size, num_items, mutation_rate, crossover_rate, generations, capacity, items):
        self.population_size = population_size
        self.num_items = num_items
        self.mutation_rate = mutation_rate
        self.crossover_rate = crossover_rate
        self.generations = generations
        self.capacity = capacity
        self.items = items
        self.population = []

    # Initialize population with random solutions (binary representation)
    def initialize_population(self):
        self.population = [[random.randint(0, 1) for _ in range(self.num_items)] for _ in range(self.population_size)]

    # Evaluate fitness of the population
    def evaluate_fitness(self):
        return [knapsack_fitness(self.items, self.capacity, individual) for individual in self.population]

    # Select individuals based on fitness (roulette wheel selection)
    def selection(self):
        fitness_values = self.evaluate_fitness()
        total_fitness = sum(fitness_values)
        if total_fitness == 0: # Avoid division by zero
            return random.choices(self.population, k=self.population_size)
        return random.choices(self.population, weights=[f / total_fitness for f in fitness_values], k=self.population_size)

    # Crossover (single-point) between two individuals
    def crossover(self, parent1, parent2):
        if random.random() < self.crossover_rate:
```

```

crossover_point = random.randint(1, self.num_items - 1)
return parent1[:crossover_point] + parent2[crossover_point:]
return parent1

# Mutation (random flip of a gene) of an individual
def mutation(self, individual):
    if random.random() < self.mutation_rate:
        mutation_point = random.randint(0, self.num_items - 1)
        individual[mutation_point] = 1 - individual[mutation_point]
    return individual

# Evolve population over generations
def evolve(self):
    self.initialize_population()
    best_solution = None
    best_fitness = 0

    for gen in range(self.generations):
        # Selection
        selected = self.selection()

        # Crossover and Mutation
        new_population = []
        for i in range(0, self.population_size, 2):
            parent1 = selected[i]
            parent2 = selected[i + 1] if i + 1 < self.population_size else selected[i]

            offspring1 = self.crossover(parent1, parent2)
            offspring2 = self.crossover(parent2, parent1)

            new_population.append(self.mutation(offspring1))
            new_population.append(self.mutation(offspring2))

        self.population = new_population

        # Evaluate fitness and track the best solution
        fitness_values = self.evaluate_fitness()
        max_fitness = max(fitness_values)
        if max_fitness > best_fitness:
            best_fitness = max_fitness
            best_solution = self.population[fitness_values.index(max_fitness)]

        print(f"Generation {gen + 1}: Best Fitness = {best_fitness}")

    return best_solution, best_fitness

# Get user input for the knapsack problem
def get_user_input():

```

```

print("Enter the number of items:")
num_items = int(input())
items = []
print("Enter the weight and value of each item (space-separated):")
for i in range(num_items):
    weight, value = map(int, input(f"Item {i + 1}: ").split())
    items.append((weight, value))

print("Enter the knapsack capacity:")
capacity = int(input())

return items, capacity, num_items

# Get user input for GEA parameters
def get_algorithm_parameters():
    print("Enter the population size:")
    population_size = int(input())
    print("Enter the mutation rate (e.g., 0.1 for 10%):")
    mutation_rate = float(input())
    print("Enter the crossover rate (e.g., 0.8 for 80%):")
    crossover_rate = float(input())
    print("Enter the number of generations:")
    generations = int(input())

    return population_size, mutation_rate, crossover_rate, generations

# Main function
if __name__ == "__main__":
    # Get user input
    items, capacity, num_items = get_user_input()
    population_size, mutation_rate, crossover_rate, generations = get_algorithm_parameters()

    # Run GEA
    gea = GeneExpressionAlgorithm(population_size, num_items, mutation_rate, crossover_rate,
                                   generations, capacity, items)
    best_solution, best_fitness = gea.evolve()

    print("\nBest Solution:", best_solution)
    print("Best Fitness (Total Value):", best_fitness)

```

```
Enter the number of items:  
5  
Enter the weight and value of each item (space-separated):  
Item 1: 10 20  
Item 2: 30 40  
Item 3: 50 60  
Item 4: 70 80  
Item 5: 90 100  
Enter the knapsack capacity:  
90  
Enter the population size:  
100  
Enter the mutation rate (e.g., 0.1 for 10%):  
0.2  
Enter the crossover rate (e.g., 0.8 for 80%):  
0.9  
Enter the number of generations:  
10  
Generation 1: Best Fitness = 120  
Generation 2: Best Fitness = 120  
Generation 3: Best Fitness = 120  
Generation 4: Best Fitness = 120  
Generation 5: Best Fitness = 120  
Generation 6: Best Fitness = 120  
Generation 7: Best Fitness = 120  
Generation 8: Best Fitness = 120  
Generation 9: Best Fitness = 120  
Generation 10: Best Fitness = 120  
  
Best solution (items selected): [1, 1, 1, 0, 0]  
Best fitness (total value): 120
```