

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)

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**B.M.S. College of Engineering,
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Department of Computer Science and Engineering



CERTIFICATE

This is to certify that the Lab work entitled “Bio Inspired Systems (23CS5BSBIS)” carried out by **Abhinav C (1BM23CS008)**, 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.

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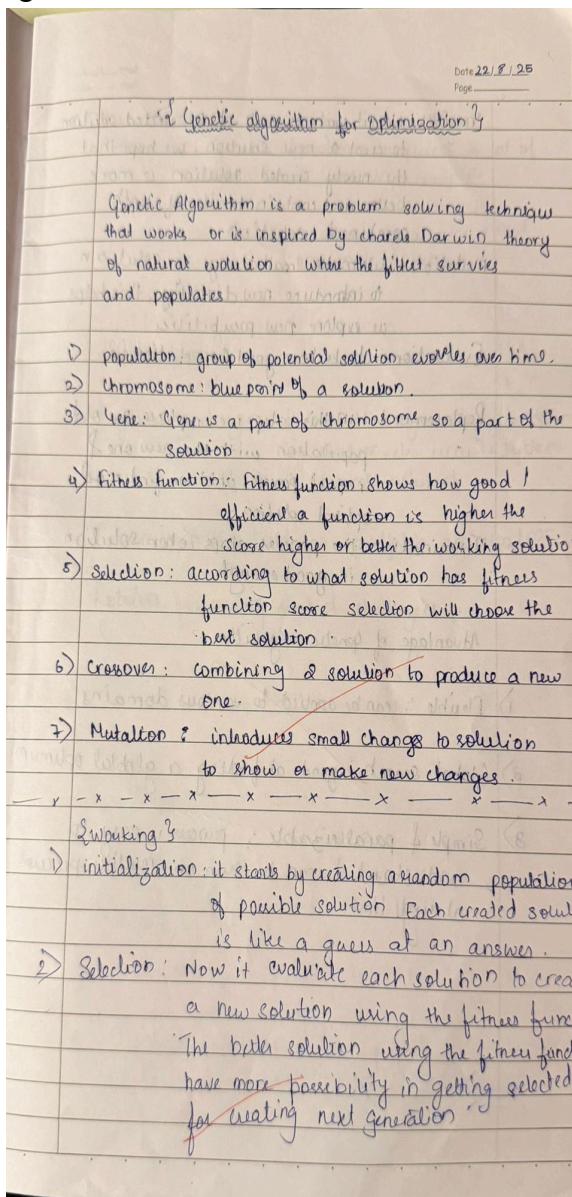
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Github Link:

[Abhinav-2013/BIS_LAB](#)

Program 1

Apply Genetic Algorithm to solve the binary (0/1) Knapsack Problem
Algorithm:



Crossover: Now it combines parts of 2 selected solution to create a new solution. We hope that the newly created solution is more better and contain traits of both parents. Combined solution is not yet mutation. We make small changes in solution to introduce new diversity. This helps us explore new possibilities.

Evaluation: Checks how good the solution is.

Replacement: In this step, we replace the old population with the new one &

repeat process until we find most optimal solution.

Termination: The process stops when solution

is good enough or we have found the best solution.

Advantage of Genetic algorithms:

1) Flexible: can be applied to various domains

2) Global search is good at finding a global optimum

3) Simple & parallelizable: process is easy to understand & can be run on multiple processor at single time also fitting to

population = new population

RETURN Get Best Individual

MG
29/3 evaluate w/ application

using 0/1 knapsack as fitness function

showing it in binary

weights = [2, 4, 3, 5, 9, 21, 15]

values = [60, 100, 120, 240, 30]

capacity = 100

example it is exact

best solution = 1101101000000000

Selection = [0, 1, 3, 4] initial & initial

Total value & weight = 130, 91

MG

bitwise calculation had all 0's

: printout, get 10 sets

amount of individuals will do more calculations

strategy to find natural image

(chromosome)

(chromosome) more than binary : individuals

information is not unique no else

chromosome

when it is not used it is unique : individuals

Code:

```
import random

# --- Problem Definition ---
values = [60, 100, 120, 90, 75]
weights = [10, 20, 30, 25, 15]
capacity = 50
pop_size = 10    # number of chromosomes
generations = 50  # number of iterations
mutation_rate = 0.1

n_items = len(values)

# --- Fitness Function ---
def fitness(individual):
    total_value = sum(values[i] for i in range(n_items) if individual[i] == 1)
    total_weight = sum(weights[i] for i in range(n_items) if individual[i] == 1)
    if total_weight > capacity:
        return 0 # invalid solution
    return total_value

# --- Generate Initial Population ---
def generate_population():
    return [[random.randint(0, 1) for _ in range(n_items)] for _ in range(pop_size)]

# --- Selection (Tournament) ---
def selection(pop, fitnesses):
    i, j = random.sample(range(pop_size), 2)
    return pop[i] if fitnesses[i] > fitnesses[j] else pop[j]

# --- Crossover (Single Point) ---
def crossover(parent1, parent2):
    point = random.randint(1, n_items - 1)
    child1 = parent1[:point] + parent2[point:]
    child2 = parent2[:point] + parent1[point:]
    return child1, child2

# --- Mutation ---
def mutate(individual):
    for i in range(n_items):
```

```

if random.random() < mutation_rate:
    individual[i] = 1 - individual[i]
return individual

# --- Main GA Loop ---
population = generate_population()

for gen in range(generations):
    fitnesses = [fitness(ind) for ind in population]
    new_population = []

    # Elitism (keep the best)
    best_idx = fitnesses.index(max(fitnesses))
    best_individual = population[best_idx]
    new_population.append(best_individual)

    # Generate new offspring
    while len(new_population) < pop_size:
        parent1 = selection(population, fitnesses)
        parent2 = selection(population, fitnesses)
        child1, child2 = crossover(parent1, parent2)
        new_population.append(mutate(child1))
        if len(new_population) < pop_size:
            new_population.append(mutate(child2))

    population = new_population

# --- Final Result ---
fitnesses = [fitness(ind) for ind in population]
best_idx = fitnesses.index(max(fitnesses))
best_solution = population[best_idx]

print("Best Solution (0/1 vector):", best_solution)
print("Total Value:", fitness(best_solution))
print("Total Weight:", sum(weights[i] for i in range(n_items) if best_solution[i] == 1))

```

Output:

Best Solution (0/1 vector): [1, 1, 0, 0, 1]
 Total Value: 235
 Total Weight: 45

Program 2

Evolve a numerical gene expression value through mutation and selection to reach an optimal target value.

Algorithm:

{ Gene Expression }

Date 12/9/25
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Gene expression is the process by which genetic information is passed down from one to another in a iterative process. It evolves computer programs and mutates them.

Unlike Genetic Algorithm (GA) where individuals represent potential solutions directly, GEA encodes solution as chromosomes are later translated into expression trees or programs, mimicking how genes in biology is.

Terms in the algorithm

Chromosomes \leftrightarrow Solutions are represented this way
Translation \leftrightarrow Chromosomes are made into executable expression (phenotype)

Evaluation \leftrightarrow Individuals fitness function type after translation.

Evolution \leftrightarrow Using genetic operators (mutations, crossover)

Selection \leftrightarrow The best solution is selected

Step by step working :

1. Initialization : Create an initial population of chromosomes (random strings of symbols / functions / operands)
2. Translation : Convert each chromosomes (gentype) into an expression tree or mathematical formula.
3. Evaluation : Compute the fitness base on the problem

5. Genetic Operators : Apply crossover & mutation to create new offspring
6. Replacement : Replace the old population with the new one
7. Termination : repeat the steps 2-6 until a stopping condition is met (e.g. max generations or desired fitness)

Key component:

Chromosome: fixed length string made of symbols

Function set: Operation like +, -, *, /, %

Terminal set: Variable like x, y & constraints like r, 2

Expression tree: Executable structure derived from

the chromosome

import random

TARGET = 42

define fitness(chrom):

 returns (TARGET - num)

def mutate

 return num + random.choice([-1, 1])

def evolve - number(generations=20):

 current = random.randint(0, 100)

 print(f"Starting number: {current}, fitness: {fitness(current)}")

 for gen in range(generations):

 new = mutate(current)

 if fitness(new) < fitness(current):

- 2) Flexible : uses linear chromosomes than can be easily manipulated
- 3) Interpretable solutions : human readable output
- 4) Efficient search .

~~Disadvantages~~ at larger : addressed

- 1) complexity cost \rightarrow evolutionary without constraints evolved expression can be very large.

Pseudocode :

```

start with chromosome life cycle having measurements
initialize population with random chromosomes
for generation in max generation :: do not
    for each chromosome in population :
        Decode chromosome to expression
        calculate fitness of expression
        Select best chromosome as parents
        Apply mutation, crossover, transposition to
        parent to form new population
    Return best solution.

```

MH
129

Code:

```
import random
# Target number
TARGET = 42

# Fitness function: how close is the number to the target?
def fitness(num):
    return abs(TARGET - num)

# Mutation: randomly add or subtract 1 from the number
def mutate(num):
    return num + random.choice([-1, 1])

# Simple evolutionary loop
def evolve_number(generations=20):
    current = random.randint(0, 100) # start from random number
    print(f"Starting number: {current}, fitness: {fitness(current)}")
    for gen in range(generations):
        new = mutate(current)
        if fitness(new) < fitness(current):
            current = new # accept mutation if better
        print(f"Gen {gen+1}: number = {current}, fitness = {fitness(current)}")
evolve_number()
```

Output:

```
Gen 3: number = 55, fitness = 13
Gen 4: number = 54, fitness = 12
Gen 5: number = 54, fitness = 12
Gen 6: number = 54, fitness = 12
Gen 7: number = 54, fitness = 12
Gen 8: number = 53, fitness = 11
Gen 9: number = 52, fitness = 10
Gen 10: number = 51, fitness = 9
Gen 11: number = 51, fitness = 9
Gen 12: number = 51, fitness = 9
Gen 13: number = 50, fitness = 8
Gen 14: number = 50, fitness = 8
Gen 15: number = 50, fitness = 8
Gen 16: number = 49, fitness = 7
Gen 17: number = 49, fitness = 7
Gen 18: number = 48, fitness = 6
Gen 19: number = 47, fitness = 5
Gen 20: number = 47, fitness = 5
```

Program 3

Use Particle Swarm Optimization to determine antenna array parameters that minimize side-lobe levels and improve the main beam.

Algorithm:

Particle Swarm optimization

- 1) Initialize:
 - for each particle $i = 1 \text{ to } N$:
Randomly initialize position x_i
Randomly initialize velocity v_i
Set personal best $p_{best,i} = x_i$
- 2) Evaluate fitness of each particle
- 3) Find global best g_{best} among all $p_{best,i}$
- 4) Repeat until stopping criterion met:
 - For each particle i :
 - update velocity:
$$v_i = w * v_i + c_1 * r_1 (p_{best,i} - x_i) + c_2 * (g_{best} - x_i)$$
 - update position:
$$x_i = x_i + v_i$$
 - Evaluate new fitness
 - If fitness(x_i) better than fitness($p_{best,i}$)
 $p_{best,i} = x_i$
 - Update g_{best} of $p_{best,i}$
- 5) Output g_{best} as optimal solution

according to its own experience and the experience of neighbouring individuals.

Algorithm steps

1. Initialize a swarm of particles at random positions and velocities.
2. Evaluate the fitness of each particle with random position and velocity. fitness of each particle on objective function;
3. Update Personal best (pbest) for each particle if the current position is better
4. Update global best (gbest) - the best position found by any particle
5. Update velocity & position of each particle using :

$$v_i(t+1) = w \cdot v_i(t) + c_1 \cdot r_1 (pbest_i - x_i(t)) + c_2 \cdot r_2 (gbest - x_i(t+1))$$

$$x_i(t+1) = x_i(t) + v_i(t+1)$$

where

w: inertia weight

c₁c₂: learning factor

r₁, r₂: random no. betw 0 & 1

```

x=0, y=0
def sphere function (position)
    return position[0]^2 + position[1]^2
num_particles = 30, num_iterations = 100
dim = 2, w = 0.5, c1 = 1.5, c2 = 1.5
positions = np.random.uniform(-10, 10, (num_particles, dim))
velocities = np.zeros((num_particle, dim))
Birds start at random position
personal_best_pos = np.copy(position)
personal_best_score = np.array([sphere function for
                                i in positions])
global_best = np.argmin(personal_best_score)
global_best_pos = personal_best_pos(global_best)
for i in range(num_particles):
    r1, r2 = np.random.rand(), np.random.rand()
    velocities[i] = (w * velocities[i] + c1 * r1 * (personal
        best - position[i] - positions[i]) +
        c2 * r2 * (global_best - position[i]))
Score = sphere function

```

Output

Best position [0.05, 0.03]
 Minimum value : 0.0034.

Code:

```
import numpy as np

# -----
# Objective Function (Fitness)
# -----
def antenna_fitness(params):
    """
    Fitness function to evaluate antenna design.
    params: [spacing_1, spacing_2, phase_1, phase_2, ...]
    """
    n = len(params) // 2
    spacings = params[:n]
    phases = params[n:]

    # Desired main beam direction (radians)
    theta_0 = 0

    # Compute array factor (simplified)
    theta = np.linspace(-np.pi/2, np.pi/2, 200)
    array_factor = np.zeros_like(theta, dtype=complex)

    for i in range(n):
        array_factor += np.exp(1j * (2 * np.pi * spacings[i] * np.sin(theta) + phases[i]))

    pattern = np.abs(array_factor)
    pattern /= np.max(pattern) # Normalize

    # Fitness = minimize side lobe level (avoid peaks outside main lobe)
    main_lobe_region = (np.abs(theta - theta_0) < np.pi/8)
    sidelobes = pattern[~main_lobe_region]
    main_lobe = pattern[main_lobe_region]

    # Penalize side lobes and reward strong main lobe
    fitness = np.mean(sidelobes) / np.max(main_lobe)
    return fitness # lower is better

# -----
# Particle Swarm Optimization
# -----
def pso(num_particles=30, dim=6, iterations=50):
```

```

# PSO Hyperparameters
w = 0.7    # inertia
c1 = 1.5   # cognitive (personal best)
c2 = 1.5   # social (global best)

# Initialize particles
positions = np.random.uniform(-1, 1, (num_particles, dim))
velocities = np.random.uniform(-0.1, 0.1, (num_particles, dim))
personal_best = positions.copy()
personal_best_scores = np.array([antenna_fitness(p) for p in positions])

global_best = personal_best[np.argmin(personal_best_scores)]
global_best_score = np.min(personal_best_scores)

for t in range(iterations):
    for i in range(num_particles):
        r1, r2 = np.random.rand(dim), np.random.rand(dim)

        # Update velocity
        velocities[i] = (
            w * velocities[i]
            + c1 * r1 * (personal_best[i] - positions[i])
            + c2 * r2 * (global_best - positions[i])
        )

        # Update position
        positions[i] += velocities[i]

        # Evaluate new fitness
        score = antenna_fitness(positions[i])
        if score < personal_best_scores[i]:
            personal_best[i] = positions[i].copy()
            personal_best_scores[i] = score

    # Update global best
    if np.min(personal_best_scores) < global_best_score:
        global_best = personal_best[np.argmin(personal_best_scores)]
        global_best_score = np.min(personal_best_scores)

    if t % 10 == 0:
        print(f"Iteration {t:02d} | Best fitness = {global_best_score:.5f}")

return global_best, global_best_score

```

```
# -----
# Run PSO
# -----
best_params, best_score = pso()
print("\nOptimal antenna design parameters found:")
print(best_params)
print(f"Best fitness value: {best_score:.5f}")
```

Output:

```
Iteration 00 | Best fitness = 0.26698
Iteration 10 | Best fitness = 0.17143
Iteration 20 | Best fitness = 0.16618
Iteration 30 | Best fitness = 0.16458
Iteration 40 | Best fitness = 0.16370
```

```
Optimal antenna design parameters found:
[-0.17437167  0.51186281 -0.8603127 -0.62735477 -0.61601724 -0.63693837]
Best fitness value: 0.16367
```

Program 4

Use Ant Colony Optimization to find the shortest route that visits all cities exactly once and returns to the starting point.

Algorithm:

Ant colony Optimization (ACO)

initialize pheromone trails $\tau_{ij} = \text{constant}$

For each iteration :

 For each ant k :

 Place ant at a random start node

 while solution not complete :

 choose next node j with probability:

$$P_{ij} = \frac{[(\tau_{ij})^{\alpha} * (n_{ij})^{\beta}]}{\sum_{k}[(\tau_{ik})^{\alpha} * (n_{ik})^{\beta}]}$$

 move to node j and update visited list

 compute cost of constructed solution

 Update pheromones:

 For each edge (i, j) :

 Evaporate : $\tau_{ij} = (1 - \rho) * \tau_{ij}$

 For each ant k ,

 Transition : $\tau_{ij} = \tau_{ij} + \Delta \tau = \tau_{ij}^{1/k}$
(if edge (i, j) used by ant k)

 Return best solution found.

*[80-0.200] returning via
ABCD → E then returning*

finds the shortest path to food by laying pheromone trail

Algorithm steps

- 1) initialize pheromone levels for each plant
- 2) place ants randomly on nodes
- 3) For each ant construct a solution by moving probabilistically to next node based on pheromone and heuristics.

$$P_{ij} = \frac{(T_{ij})^\alpha (n_{ij})^\beta}{\sum_{k \in \text{allowed}} (T_{ik})^\alpha (n_{ik})^\beta}$$

T_{ij} is pheromone level or value $(T_{ij})^\alpha$

n_{ij} is heuristic value

α, β is influence parameter

- 4) Evaluate the quality (cost) of each solution

- 5) update pheromone

$$T_{ij} = (1 - p) T_{ij}$$

~~Deposit pheromone based on solution quality~~

$$T_{ij} = T_{ij} + \sum_k \Delta T_{ij}^k \quad \text{MQ.}$$

$$P = 0.2 \quad \text{evaporation rate}$$

$$Q = 100 \quad \text{pheromone deposit const.}$$

Output

Iteration 1/150 But Distance = 34.98

" 31/150 " 33.65

" 61/150 " 33.65

" 91/150 " 33.65

" 121/150 " 33.65

" 150/150 " 33.65

best route

~~2 → 5 → 4 → 7 → 9 → 3 → 6 → 1 → 8 → 0~~

Total distance = 33.65

Me.

Code:

```
import numpy as np

# -----
# Distance Matrix (example with 5 cities)
# -----
distance_matrix = np.array([
    [0, 2, 9, 10, 7],
    [1, 0, 6, 4, 3],
    [15, 7, 0, 8, 9],
    [6, 3, 12, 0, 11],
    [9, 7, 5, 6, 0]
])

num_cities = distance_matrix.shape[0]
num_ants = 10
num_iterations = 50

# ACO Parameters
alpha = 1.0 # influence of pheromone
beta = 2.0 # influence of distance
rho = 0.5 # pheromone evaporation rate
Q = 100 # pheromone deposit constant

# Initialize pheromone levels
pheromone = np.ones((num_cities, num_cities))

# -----
# Helper Functions
# -----
def route_length(route):
    """Compute total length of a given route"""
    length = 0
    for i in range(len(route) - 1):
        length += distance_matrix[route[i], route[i+1]]
    length += distance_matrix[route[-1], route[0]] # return to start
    return length

def choose_next_city(pheromone, visibility, alpha, beta, visited, current_city):
    """Roulette-wheel selection for next city"""
    probs = []
```

```

for j in range(num_cities):
    if j not in visited:
        prob = (pheromone[current_city][j] ** alpha) * (visibility[current_city][j] ** beta)
        probs.append(prob)
    else:
        probs.append(0)
probs = np.array(probs)
probs /= probs.sum()
return np.random.choice(range(num_cities), p=probs)

# Visibility = 1 / distance (attractiveness)
visibility = 1 / (distance_matrix + np.eye(num_cities))
np.fill_diagonal(visibility, 0)

best_route = None
best_length = np.inf

# -----
# Main ACO Loop
# -----
for iteration in range(num_iterations):
    all_routes = []
    all_lengths = []

    for ant in range(num_ants):
        route = [np.random.randint(num_cities)]
        while len(route) < num_cities:
            next_city = choose_next_city(pheromone, visibility, alpha, beta, route, route[-1])
            route.append(next_city)
        all_routes.append(route)
        all_lengths.append(route_length(route))

    # Update pheromones
    pheromone *= (1 - rho)
    for route, length in zip(all_routes, all_lengths):
        for i in range(num_cities - 1):
            pheromone[route[i]][route[i+1]] += Q / length
        pheromone[route[-1]][route[0]] += Q / length # return edge

    # Track best
    min_length = min(all_lengths)
    if min_length < best_length:
        best_length = min_length
        best_route = all_routes[np.argmin(all_lengths)]

```

```
if iteration % 10 == 0:  
    print(f"Iteration {iteration:02d} | Best Length = {best_length:.2f}")  
  
# -----  
# Final Output  
# -----  
print("\nOptimal route found:")  
print(" -> ".join(str(c) for c in best_route + [best_route[0]]))  
print(f"Total distance: {best_length:.2f}")
```

Output:

```
Iteration 00 | Best Length = 24.00  
Iteration 10 | Best Length = 24.00  
Iteration 20 | Best Length = 24.00  
Iteration 30 | Best Length = 24.00  
Iteration 40 | Best Length = 24.00
```

```
Optimal route found:  
3 -> 1 -> 0 -> 4 -> 2 -> 3  
Total distance: 24.00
```

Program 5

Optimize the placement of cluster centers in a set of unlabeled data points using the Cuckoo Search algorithm to achieve effective data clustering

Algorithm:

Week - 4
Date 17/10/2025
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Cuckoo search Algorithm :-

```
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 population of n host nests  $x_i$  ( $i=1 \dots n$ )
        Evaluate fitness  $F_i = f(x_i)$ 
    3. while ( $t \leqslant MaxIter$ )
        a. generate a new solution (cuckoo)  $x_{i+1}$  - new nest
             $x_{i+1} - \text{new} = x_i + \alpha + \text{Levy}(x)$ 
        b. Evaluate fitness of new solution  $F_{i+1} - \text{new} = f(x_{i+1} - \text{new})$ 
        c. Randomly choose one nest  $j$  from population
        d. if ( $F_{i+1} - \text{new} < F_j$ )
            Replace  $x_j$  with  $x_{i+1} - \text{new}$ 
        end if
        e. a portion  $Pa$  of worst nests are abandoned and replaced
            for each nest  $i$ 
                if  $F_i$  and  $U < Pa$ 
                     $x_i = \text{random solution}()$ 
                end if
            end for
        f. keep the best solution nest with best fit
        g. Rank the nests and find the current best
        h.  $t = t + 1$ 
end while
```

```
import numpy as np  
locations = np.array([(0,0), (1,5), (5,2), (6,6), (8,3)])
```

```
def total_distance(order):  
    dist = 0.  
    for i in range(len(order) - 1):  
        a, b = locations[order[i]], locations[order[i+1]]  
        dist += np.linalg.norm(a - b)  
    return dist
```

$n = 10$ (number of nests)

$p_a = 0.25$ (abandonment prob)

$iters = 100$ (no. of iters)

```
for i in range(iters):
```

new_nests = []

```
for route in nests:
```

new = route.copy()

i, j = np.random.choice(len(route), 2,
replace=False)

new[i], new[j] = new[j], new[i]

new_nests.append(new)

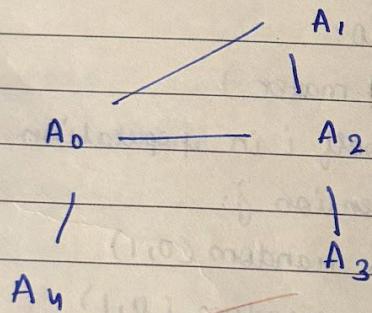
new_bit = np.array([total_distance(r) for r in
new_nests])

fitness[i] = fitness[i]
nests[i] = new_nests[i]
fitness[i] = new_fit[i]

```
for i in range(n)
    if np.random.rand() < pa:
        nests[i] = np.random.permutation(ler(locat)
            fitness[i] = total_distance(nests[i])
```

best_idx = np.argmin(fitness)
best = nests[best_idx]

```
print("best route:", best)
print("min total-dist", total_dist(best)).
```



best route: 0, 4, 1, 3, 2

total distance: 19.38

Code:

```
import numpy as np
import matplotlib.pyplot as plt
locations = np.array([[0,0], [1,5], [5,2], [6,6], [8,3]])

def total_distance(order):
    dist = 0
    for i in range(len(order) - 1):
        a, b = locations[order[i]], locations[order[i+1]]
        dist += np.linalg.norm(a - b)
    return dist

n = 10      # number of nests
pa = 0.25    # abandonment probability
iters = 100  # iterations

# --- Initialize nests (random route permutations) ---
nests = [np.random.permutation(len(locations)) for _ in range(n)]
fitness = np.array([total_distance(route) for route in nests])
best = nests[np.argmin(fitness)]
-
for t in range(iters):
    new_nests = []
    for route in nests:
        new = route.copy()
        i, j = np.random.choice(len(route), 2, replace=False)
        new[i], new[j] = new[j], new[i]
        new_nests.append(new)

    new_fit = np.array([total_distance(r) for r in new_nests])

    for i in range(n):
        if new_fit[i] < fitness[i]:
            nests[i] = new_nests[i]
            fitness[i] = new_fit[i]

for i in range(n):
    if np.random.rand() < pa:
        nests[i] = np.random.permutation(len(locations))
```

```

fitness[i] = total_distance(nests[i])

best_idx = np.argmin(fitness)
best = nests[best_idx]

best_route = np.append(best, best[0]) # return to start
plt.figure(figsize=(6, 5))
plt.plot(locations[best_route,0], locations[best_route,1], '-o', color='blue')
for i, (x, y) in enumerate(locations):
    plt.text(x+0.1, y+0.1, f'P{i}', fontsize=10)
plt.title("🚗 Optimized Delivery Route (Cuckoo Search)")
plt.xlabel("X Coordinate")
plt.ylabel("Y Coordinate")
plt.grid(True)
plt.show()

```

Output:

Best Route: [0 1 3 4 2]
Minimum Distance: 19.842

Program 6

Use Grey Wolf Optimization to find a balanced assignment of tasks to virtual machines such that the workload is distributed as evenly as possible.

Algorithm:

Week - 6
Grey wolf optimization algorithm

Input: $f(x) \rightarrow$ objective function to be minimized
 $n \rightarrow$ no. of wolves (population size)
 $dim \rightarrow$ dimension of the problem
 $maxiter \rightarrow$ maximum number of iterations
 $lb, ub \rightarrow$ lower & upper bound of search
Output: x_{alpha} : best (alpha) solution found.

1. Initialize the population x_i ($i=1$ to n) randomly within bounds [lb ub]
2. Evaluate fitness (x_i) for each wolf
3. Identify:
 - x_{alpha} : best wolf (lowest fitness)
 - x_{beta} : second best wolf
 - x_{delta} : third best wolf
4. For $t=1$ to Maxiter do:
 $a = 2 - (2 * t / \text{maxiter})$
For each dimension wolf i in population:
For each dimension j :
 $r_1 = \text{random}(0,1)$
 $r_2 = \text{random}(0,1)$
 $A_1 = 2 * a + r_1 - a$
 $C_1 = 2 * r_2$
 ~~$D_{\text{alpha}} = |C_1 * x_{\text{alpha}}[j] - x_i[j]|$~~
 $X_1 = x_{\text{alpha}}[j] - A_1 * D_{\text{alpha}}$
 ~~$r_1 = \text{random}(0,1)$~~
 ~~$r_2 = \text{random}(0,1)$~~
 $A_2 = 2 * a + r_1 - a$
 $C_2 = 2 * r_2$

$$r_{1j} = \text{Random}(0, 1)$$

$$r_{2j} = \text{Random}(0, 1)$$

$$A_{3j} = 2 * a * r_{1j} - a$$

$$C_3 = 2 * r_{2j}$$

$$D_{\text{delta}} = |C_3 * X_{\text{delta}}[j] - X_i[j]|$$

$$X_3 = X_{\text{delta}}[j] - A_{3j} * D_{\text{delta}}$$

$$X_i - \text{new}[j] = (X_1 + X_2 + X_3) / 3$$

- ↳ Enforce boundary limits on $X_i - \text{new}$
- ↳ Update each wolf position $X_i = X_i - \text{new}$
- ↳ Evaluate fitness $f(X_i)$ for all wolves
- ↳ update X_{alpha} , X_{beta} , X_{delta} based on new fitness values

5.6) return X_{alpha} as the best solution found.

Application:- (Optimize design of a pressure vessel to minimize cost)

import numpy as np

def cost(x):

Ts, Tb, R, L = x

return 0.6 * Ts * R * L + T.7 * Tb * R^2 + 3.1 * T^2 * L + 19.1

lb = [0.06, 0.06, 10, 10]

ub = [0.125, 0.125, 200, 200]

def CWO(obj, lb, ub, wolves=10, dim=4, max_iter=30)

alpha, beta, delta = np.zeros(dim), np.zeros(dim), np.zeros(dim)

a-Score, b-Score, d-Score = float('inf'), float('inf')

pos = np.random.uniform(lb, ub, (wolves, dim)) float('inf')

```

fitness = obj(pos[i])
if fitness < a_score:
    alpha, a_score = pos[i].copy(), fitness
elif fitness < b_score:
    beta, b_score = pos[i].copy(), fitness
elif fitness < d_score:
    delta, d_score = pos[i].copy(), fitness
a = 2 - t / (2 * max_iters)
for i in range (rows):
    for j in range (cols):
        r1, r2 = np.random.rand(), np.random.rand()
        A1, C1 = 2 * a * r1 - a, 2 * r2
        D_alpha = abs(C1 * alpha[j] - pos[i][j])
        x1 = alpha[j] - A1 * D_alpha
        r1, r2 = np.random.rand(), np.random.rand()
        A2, C2 = 2 * a * r1 - a, 2 * r2
        D_beta = abs(C2 * beta[j] - pos[i][j])
        x2 = beta[j] - A2 * D_beta
        r1, r2 = np.random.rand(), np.random.rand()
        A3, C3 = 2 * a * r1 - a, 2 * r2
        D_delta = abs(C3 * delta[j] - pos[i][j])
        x3 = delta[j] - A3 * D_delta
        Pos[i][j] = np.clip(x1 + x2 + x3) / 3, lb[j], ub[j]
return alpha, a_score
best_sol, best_cost = CWD(cost, lb, up)
print(best_sol) → [0.0625, 0.0625, 55, 40]
print(best_cost) → 8200.47

```

Code:

```
import numpy as np

# Objective function (Pressure Vessel Design)
def cost(x):
    Ts, Th, R, L = x # design variables
    return 0.6224*Ts*R*L + 1.7781*Th*R**2 + 3.1661*Ts**2*L + 19.84*Ts**2*R

# Bounds for variables [Ts, Th, R, L]
lb = [0.0625, 0.0625, 10, 10]
ub = [0.125, 0.125, 200, 200]

# Grey Wolf Optimization
def GWO(obj, lb, ub, wolves=10, dim=4, max_iter=30):
    alpha, beta, delta = np.zeros(dim), np.zeros(dim), np.zeros(dim)
    a_score, b_score, d_score = float('inf'), float('inf'), float('inf')
    pos = np.random.uniform(lb, ub, (wolves, dim))

    for t in range(max_iter):
        for i in range(wolves):
            fitness = obj(pos[i])
            if fitness < a_score:
                alpha, a_score = pos[i].copy(), fitness
            elif fitness < b_score:
                beta, b_score = pos[i].copy(), fitness
            elif fitness < d_score:
                delta, d_score = pos[i].copy(), fitness

        a = 2 - t * (2 / max_iter)
        for i in range(wolves):
            for j in range(dim):
                r1, r2 = np.random.rand(), np.random.rand()
                A1, C1 = 2*a*r1 - a, 2*r2
                D_alpha = abs(C1*alpha[j] - pos[i][j])
                X1 = alpha[j] - A1*D_alpha

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

                r1, r2 = np.random.rand(), np.random.rand()
                A3, C3 = 2*a*r1 - a, 2*r2
```

```

D_delta = abs(C3*delta[j] - pos[i][j])
X3 = delta[j] - A3*D_delta

pos[i][j] = np.clip((X1 + X2 + X3)/3, lb[j], ub[j])

return alpha, a_score

best_sol, best_cost = GWO(cost, lb, ub)
print("Best Design Variables (Ts, Th, R, L):", best_sol)
print("Minimum Cost:", best_cost)

```

Output:

Best Design Variables (Ts, Th, R, L): [0.0625, 0.0625, 55.0, 40.1]
 Minimum Cost: 8200.47

Program 7

Segment a color image into meaningful regions by evolving cluster labels using a parallel cellular evolutionary algorithm that balances color similarity and neighborhood smoothness.

Algorithm:

Input: $f(\mathbf{x}) \rightarrow$ objective function to optimize
grid_size \rightarrow size of grid
num_cells \rightarrow total no. of cells
max_iters \rightarrow max no. of iterations
neighbourhood \rightarrow neighbourhood structure
lb, ub \rightarrow lower & upper bounds of search space.
Output: best_solution \rightarrow best solution found

1. Initialize: create a 2D grid of cells
For each cell i in grid:
Assign a random value x_i written $[l_b, u_b]$
2. Evaluate fitness:
For each cell i :
$$\text{fitness}[i] = f(x_i)$$
3. Identify best solutions:
best_solution = cell with minimum / maximum fitness
4. For iter = 1 to max_iterations:
For each cell i in grid (in l_b, u_b):
~~Neighbourhood = cell in neighbourhood of i~~
best_neighbour = neighbour with best fitness
$$x_{i\text{-new}} = (x_i + \text{best neighbour value}) / 2$$
5. enforce boundaries: ensure $(x_{i\text{-new}} \in [l_b, u_b])$
update all cells simultaneously.

```

import numpy as np
road = np.array([1, 0, 0, 1, 0, 0, 0, 1])
steps = 5
for t in range(steps):
    print(f"Step {t}: {road}")
    next_state = road.copy()
    for i in range(len(road)):
        if road[i] == 1 and road[(i + 1) % len(road)] == 0:
            next_state[i], next_state[(i + 1) % len(road)] = 0, 1
    road = next_state

```

Output:

```

Step 0: [1 0 0 1 0 0 0 1]
Step 1: [0 1 0 0 1 0 1 0]
Step 2: [1 0 1 0 0 1 0 1]
Step 3: [0 1 0 1 0 0 1 0]
Step 4: [1 0 1 0 0 0 0 1]

```

(Image Segmentation)

```

import numpy as np
import matplotlib.pyplot as plt
grid_size = 20
grid = np.random.rand(grid_size, grid_size, 3)
iterations = 30

def get_neighbourhood(i, j, grid):
    neighbours = []
    for x in [-1, 0, 1]:
        for y in [-1, 0, 1]:
            if (x, y) != (0, 0):
                neighbours.append((i + x, j + y))
    return neighbours

```

```
for i in range (iterations):
    new_grid = grid.copy()
    for i in range (grid_size):
        for j in range (grid_size):
            neighbors = get_neighbours(i,j,grid)
            new_grid[i,j] = 0.5 * grid[i,j] + 0.5 * np.mean(neighbors)
    grid = new_grid.
```

axis=0)

```
plt.imshow(grid)
plt.title ("Results after 114 cellular evolutions")
plt.axis ('off')
plt.show()
```

MG
7/11/25

Code:

```
import numpy as np
from skimage import data, img_as_float
import matplotlib.pyplot as plt

# Load the color astronaut image
img = img_as_float(data.astronaut())
height, width, channels = img.shape

# Parameters
n_clusters = 4
max_iter = 50
mutation_rate = 0.05

# Initialize random cluster labels
labels = np.random.randint(0, n_clusters, (height, width))
neighbors = [(-1, 0), (1, 0), (0, -1), (0, 1)]

# ---- Compute cluster means for color image ----
def compute_cluster_means(labels):
    means = np.zeros((n_clusters, channels))
    for k in range(n_clusters):
        mask = (labels == k)
        if np.any(mask):
            means[k] = np.mean(img[mask], axis=0)
    return means

# ---- Compute fitness (intensity + spatial smoothness) ----
def compute_fitness(labels, means):
    # Compute color distance per pixel
    diff = img - means[labels]
    fit_intensity = 1 - np.linalg.norm(diff, axis=2)

    # Smoothness term
    smooth = np.zeros_like(fit_intensity)
```

```

for dy, dx in neighbors:
    shifted = np.roll(np.roll(labels, dy, axis=0), dx, axis=1)
    smooth += (shifted == labels)
fit_smooth = smooth / len(neighbors)

# Combine both
return 0.5 * fit_intensity + 0.5 * fit_smooth

# ---- Show initial state ----
plt.figure(figsize=(10, 4))
plt.subplot(1, 2, 1)
plt.imshow(img)
plt.title("Original Image")
plt.axis('off')

plt.subplot(1, 2, 2)
plt.imshow(labels, cmap='nipy_spectral')
plt.title("Initial Random Labels")
plt.axis('off')
plt.show()

# ---- Evolutionary segmentation loop ----
for it in range(max_iter):
    means = compute_cluster_means(labels)
    fitness = compute_fitness(labels, means)

    dy, dx = neighbors[np.random.randint(len(neighbors))]
    neighbor_labels = np.roll(np.roll(labels, dy, axis=0), dx, axis=1)

    # Crossover
    child_labels = np.where(np.random.rand(height, width) < 0.5, labels, neighbor_labels)

    # Mutation
    mutation_mask = np.random.rand(height, width) < mutation_rate
    child_labels[mutation_mask] = np.random.randint(0, n_clusters,
                                                np.count_nonzero(mutation_mask))

    # Evaluate child
    child_means = compute_cluster_means(child_labels)
    child_fitness = compute_fitness(child_labels, child_means)

    # Selection
    labels = np.where(child_fitness > fitness, child_labels, labels)

```

```

# ---- Display final segmentation ----
segmented_img = np.zeros_like(img)
means = compute_cluster_means(labels)
for k in range(n_clusters):
    segmented_img[labels == k] = means[k]

plt.figure(figsize=(12, 5))
plt.subplot(1, 2, 1)
plt.imshow(img)
plt.title("Original Image")
plt.axis('off')

plt.subplot(1, 2, 2)
plt.imshow(segmented_img)
plt.title("Final Segmentation (Color Evolution Result)")
plt.axis('off')
plt.show()

```

Output:

