
VISVESVARAYA TECHNOLOGICAL UNIVERSITY

“JnanaSangama”, Belgaum -590014, Karnataka.



LAB RECORD

Bio Inspired Systems (23CS5BSBIS)

Submitted by **Sanchit Kashyap(1BM23CS298)**

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,
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 **Sanchit Kashyap(1BM23CS298)**, who is bona fide 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.

Rohith Vaidya K Assistant Professor Department of CSE, BMSCE	Dr. Selva Kumar Professor & HOD Department of CSE, BMSCE
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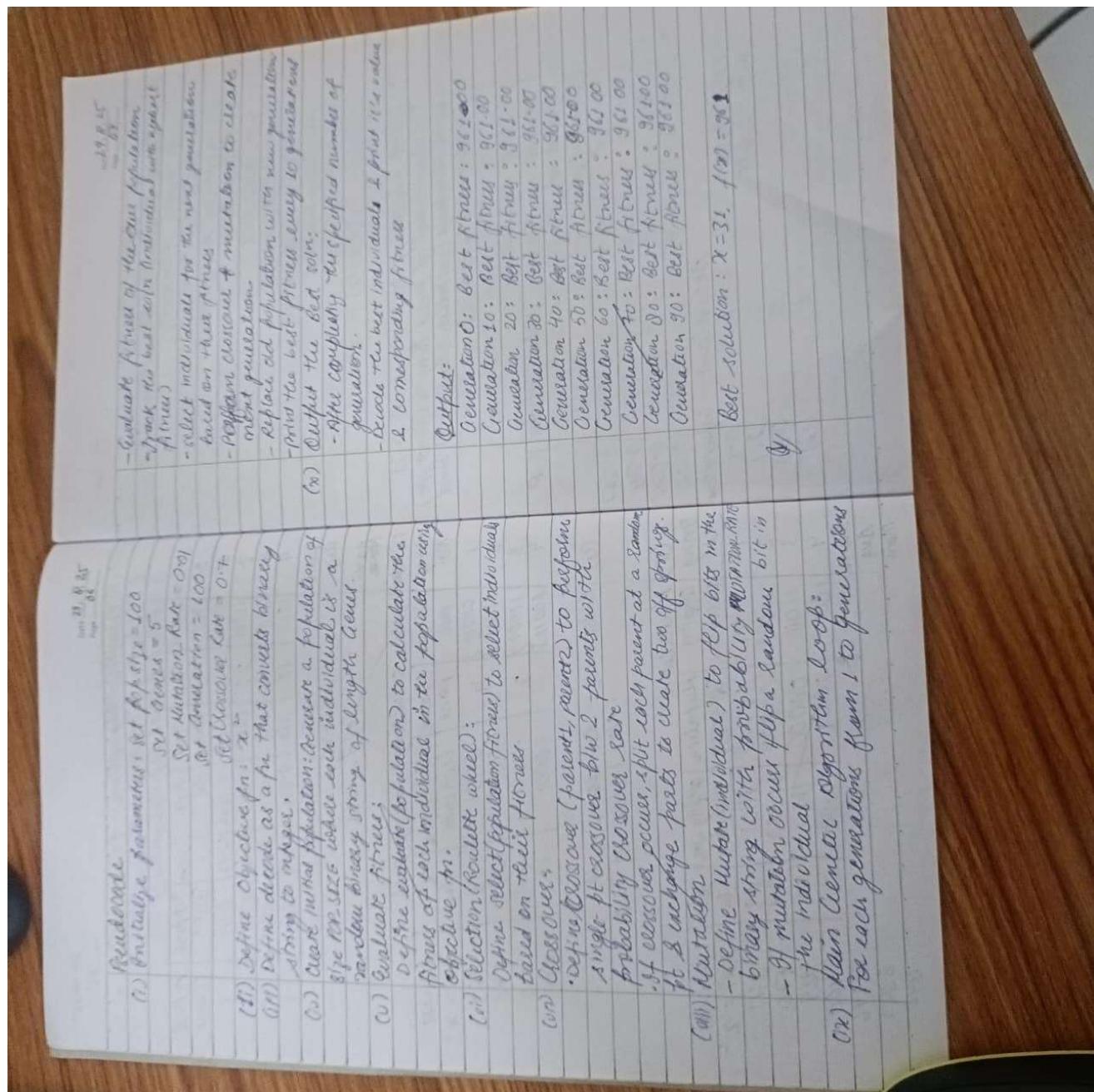
Github Link:

https://github.com/sanchit901/BIS_LAB_1BM23CS298

Program 1

Genetic Algorithm for Optimization Problems

Algorithm:



Code:

```
import numpy as np  
  
# Objective function to maximize
```

```

def fitness_function(x):
    return x**2

# Initialize parameters
population_size = 50 mutation_rate
= 0.1 crossover_rate = 0.7
num_generations = 50 lower_bound =
-10 upper_bound = 10

# Create initial population def
initialize_population(size, lower, upper):      return
np.random.uniform(lower, upper, size)

# Evaluate fitness for the population def
evaluate_fitness(population):
    return np.array([fitness_function(x) for x in population])
# Selection using roulette wheel selection def
select_parents(population, fitness):
    total_fitness = np.sum(fitness)      selection_probs = fitness /
total_fitness      parents_indices =
np.random.choice(len(population), size=2, p=selection_probs)
return population[parents_indices]

# Crossover to create offspring def
crossover(parent1, parent2):
    if np.random.rand() < crossover_rate:
        return (parent1 + parent2) / 2 # Linear crossover      return
parent1

# Mutation to introduce diversity def
mutate(offspring):
    if np.random.rand() < mutation_rate:
        return np.random.uniform(lower_bound, upper_bound)      return
offspring

# Genetic Algorithm main function def genetic_algorithm():      #
Initialize population      population =
initialize_population(population_size, lower_bound, upper_bound)
    for generation in range(num_generations):

```

```

        # Evaluate fitness of the population
fitness = evaluate_fitness(population)

        # Track the best solution
best_fitness_idx = np.argmax(fitness)
best_solution = population[best_fitness_idx]
best_fitness_value = fitness[best_fitness_idx]

        print(f"Generation {generation}: Best Solution = {best_solution},
Fitness = {best_fitness_value}")

        # Create the next generation
new_population = []           for _ in
range(population_size):
    parent1, parent2 = select_parents(population,
fitness)                 offspring = crossover(parent1, parent2)
offspring = mutate(offspring)
new_population.append(offspring)
population =
np.array(new_population)

        # Final evaluation      final_fitness =
evaluate_fitness(population)    best_fitness_idx =
np.argmax(final_fitness)       best_solution =
population[best_fitness_idx]   best_fitness_value =
final_fitness[best_fitness_idx]
    return best_solution,
best_fitness_value

# Run the genetic algorithm
best_solution, best_fitness_value = genetic_algorithm()
print(f"Best Solution Found: x = {best_solution}, f(x) =
{best_fitness_value}")

```

Output :

```

Generation 0: Best Solution = -9.967365011554792, Fitness = 99.34836527356666
Generation 1: Best Solution = -9.169251894044368, Fitness = 84.07518029643623
Generation 49: Best Solution = 9.123059138454053, Fitness = 83.23020804373002
Best Solution Found: x = 9.05670095588789, f(x) = 82.02383220438064

```

Program 2

Particle Swarm Optimization for Function Optimization

Algorithm:

→ Particle Swarm Optimization (PSO)
Pseudo code:-

• $P = \text{particle initialization}$
for $t=1 \rightarrow \text{more}$
for each particle p in P do
 $f_p = f(p)$

if f_p is better than $f(p_{best})$
 $p_{best} = p$
end for
end for

$g_{best} = \text{best from } P$
for each particle p in P do
 $v_i^{t+1} = v_i^t + C_1 U_i^t (p_{best} - p_i^t) + C_2 U_s^t (g_{best} - p_i^t)$
intra personal social
influence influence
 $p_i^{t+1} = v_i^{t+1} + U_i^{t+1}$

Code:

```
import numpy as np

# Step 1: Define the Problem def
fitness_function(position):
    # Example: Minimize the Sphere function      return
np.sum(position**2)

# Step 2: Initialize Parameters def
initialize_parameters():
    params = {
        'N': 50,          # Number of particles
        'dim': 2,         # Dimensionality of the problem
        'max_iter': 200,  # Maximum number of iterations
        'minx': -10,     # Minimum bound for position
        'maxx': 10,       # Maximum bound for position
        'w': 0.5,         # Inertia weight
        'c1': 1.5,        # Cognitive coefficient
        'c2': 1.5         # Social coefficient
    }      return params

# Step 3: Initialize Particles class Particle:    def
__init__(self, position, velocity):
    self.position = position      self.velocity
= velocity      self.bestPos = position.copy()
self.bestFitness = float('inf')
    def initialize_swarm(N, dim, minx, maxx):
        swarm = []      for _ in
range(N):
            position = np.random.uniform(minx, maxx, dim)
velocity = np.random.uniform(-1, 1, dim)
swarm.append(Particle(position, velocity))      return swarm

# Step 4: Evaluate Fitness def
evaluate_fitness(swarm):      for
particle in swarm:
    particle.fitness = fitness_function(particle.position)
# Step 5: Update Velocities and Positions def update_particles(swarm,
best_pos_swarm, w, c1, c2, minx, maxx):
```

```

        for particle in swarm:
            r1, r2 = np.random.rand(), np.random.rand()           particle.velocity
= (w * particle.velocity +                                r1 * c1 * (particle.bestPos
- particle.position)                                r2 * c2 * (best_pos_swarm -
particle.position))          particle.position += particle.velocity      #
Clip position to be within bounds                  particle.position =
np.clip(particle.position, minx, maxx)
# Step 6: Iterate
def pso():
    params = initialize_parameters()      swarm =
initialize_swarm(params['N'], params['dim'], params['minx'],
params['maxx'])      best_pos_swarm = swarm[0].position.copy()
best_fitness_swarm = float('inf')
    for _ in
range(params['max_iter']):
        evaluate_fitness(swarm)
        for particle in swarm:             if
particle.fitness < particle.bestFitness:
particle.bestFitness = particle.fitness
particle.bestPos = particle.position.copy()           if
particle.fitness < best_fitness_swarm:
best_fitness_swarm = particle.fitness
best_pos_swarm = particle.position.copy()
        update_particles(swarm,       best_pos_swarm,       params['w'],
params['c1'], params['c2'], params['minx'], params['maxx'])

    # Step 7: Output the Best Solution
return best_pos_swarm, best_fitness_swarm
best_position, best_fitness =
pso()

print("Best Position:", best_position) print("Best
Fitness:", best_fitness)

Output :

```

```

Best Position: [-9.19971249e-25  1.71937901e-24]
Best Fitness: 3.802611270068504e-48

```

Program 3

Ant Colony Optimization for the Traveling Salesman Problem

Algorithm:

Date 10/10/25
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→ Ant Colony Optimization for the Traveling Salesman Problem

Posn	Fitness value
1	2
-	2
-	0.5
-	2
-	0.125

Fitness value
 2
 2
 0.5
 2
 0.125

Pseudo-code:-
 Input:
 - cities: list of city coordinates
 n_ants: no. of ants
 n_iterations: no. of iterations
 alpha: pheromone importance factor
 beta: heuristic imp. factor (inverse distance)
 rho: pheromone evaporation rate
 initial_pheromone: initial pheromone value on edges

Output:
 best_tour: seq. of cities representing the shortest route found
 best_length: length of the best tour

Procedure:
 1. Calc dist matrix b/w all pairs of cities
 2. Initialize pheromone matrix with initial pheromone for all edges
 3. Calc heuristic matrix as inverse of dist. mat.
 4. Initialize best_length to a large no. and best_tour as empty.
 5. For iteration = 1 to n_iterations do:
 a. For each ant = 1 to n_ants do:
 i. Randomly select a start city for the ant
 ii. Initialize visited cities with start city
 iii. While there are unvisited cities do:
 - for each unvisited city, calc prob. proportionally:

$$\frac{\text{pheromone}[\text{current city}]^{\alpha} \cdot \text{alpha}}{\text{heuristic}[\text{current city}]^{\beta} \cdot \text{beta}}$$

- Normalize the probabilities
- Select the next city based on the prob distribution
- Add the selected city to visited cities
- iv. Calc the length of the completed tour (including return to start).

5. After all ants have completed tours:

i. Evaporate pheromone on all edges:

$$\text{pheromone}_{i,j} = (1 - rho) * \text{pheromone}_{i,j}$$

ii. For each ant:

- deposit pheromone on edges in the ant's tour proportionality to $1/\text{tour length}$:

$$\text{pheromone}_{\text{city } i}[\text{city } j] += Q / \text{tour length}$$

6. Update best tour and best length if any ant found a better tour.

7. Return best tour and best length

Output:-

$$\text{cities} = (0, 0), (1, 5), (5, 2), (6, 6), (8, 3), (7, 7)$$

~~Routes = 20~~

$$n_ants = 20$$

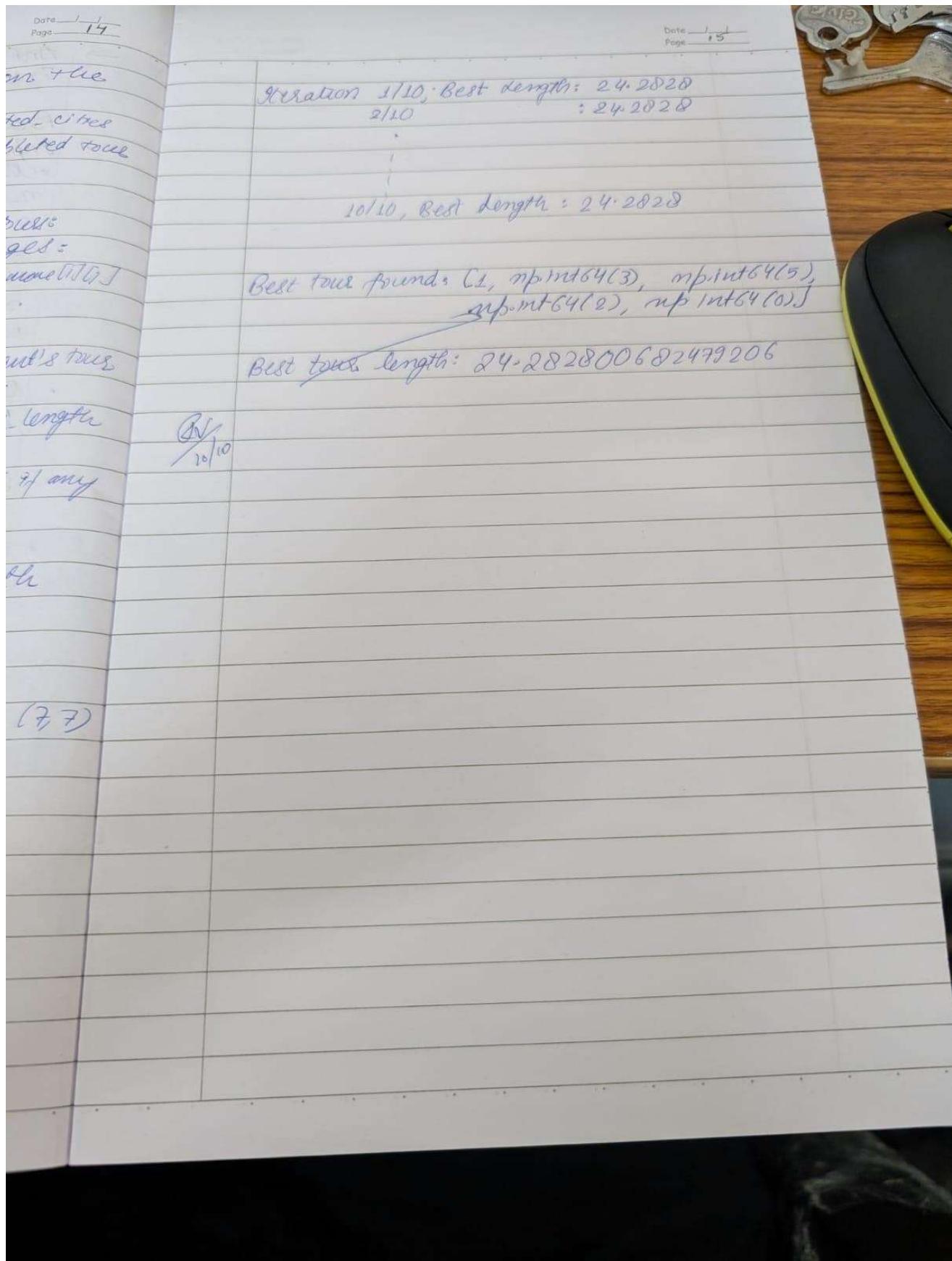
$$n_iterations = 100$$

$$\alpha = 1$$

$$\beta = 5$$

$$\rhoho = 0.5$$

$$\text{initial_pheromone} = 1$$



Code:

```
import numpy as np
random
class
AntColony:
    def __init__(self, cities, num_ants=10, alpha=1.0, beta=2.0, rho=0.5,
iterations=100):
        self.cities = cities          self.num_ants = num_ants
self.alpha = alpha          self.beta = beta          self.rho = rho
self.iterations = iterations      self.num_cities = len(cities)
self.pheromone = np.ones((self.num_cities, self.num_cities))
self.distance = self.calculate_distances()      def
calculate_distances(self):
    distances = np.zeros((self.num_cities, self.num_cities))
for i in range(self.num_cities):          for j in range(i + 1,
self.num_cities):          distances[i][j] =
distances[j][i] = np.linalg.norm(np.array(self.cities[i]) -
np.array(self.cities[j]))      return distances      def
select_next_city(self, current_city, visited):
    probabilities = []          for next_city
in range(self.num_cities):          if
next_city not in visited:
        pheromone = self.pheromone[current_city][next_city]  **
self.alpha          heuristic = (1 /
self.distance[current_city][next_city])  **          self.beta
probabilities.append(pheromone * heuristic)          else:
        probabilities.append(0)
total = sum(probabilities)
probabilities = [p / total for p in probabilities]
return np.random.choice(range(self.num_cities), p=probabilities)
def construct_solution(self):
    for _ in range(self.num_ants):
        visited = [0]
current_city = 0          for _ in range(1,
self.num_cities):
        next_city = self.select_next_city(current_city, visited)
```

```

        visited.append(next_city)
current_city = next_city           visited.append(0)  #
Return to starting city           yield visited      def
update_pheromones(self, solutions):
    self.pheromone *= (1 - self.rho) # Evaporation      for
solution in solutions:
    length = self.calculate_tour_length(solution)
pheromone_deposit = 1 / length          for i in
range(len(solution) - 1):
    self.pheromone[solution[i]][solution[i + 1]] +=
pheromone_deposit      def calculate_tour_length(self, solution):
    return sum(self.distance[solution[i]][solution[i + 1]] for i in
range(len(solution) - 1))
def run(self):
    best_solution = None
best_length = float('inf')          for _ in
range(self.iterations):
    solutions = list(self.construct_solution())
self.update_pheromones(solutions)      for solution
in solutions:
    length = self.calculate_tour_length(solution)
if length < best_length:            best_length =
length                                best_solution = solution
return best_solution, best_length cities = [(0, 0), (1, 2), (2,
1), (4, 4), (2, 4)] aco = AntColony(cities)
best_route, best_distance = aco.run()
print("Best Route:", best_route)
print("Best Distance:", best_distance)

```

Output :

```

Best Route: [0, 1, 4, 3, 2, 0]
Best Distance: 12.313755207963359

```

Program 4 Cuckoo
Search (CS)

Algorithm:

→ Cuckoo Search
Brute Force:-

Cuckoo Search:-

Initialize population of m best nests (solutions)

$\mathcal{N}(x^1, x^2, \dots, x^m)$

Define objective fn $f(x) = (x_1, x_2, \dots, x_d)^T$

And find our best x^{best} α -best among the nests

While $t < M$ iterations OR stopping criteria

for each nest $i \in \mathcal{N}_t$, do

Generate a new x_i^{new} at x_i^{old} by Levy flight

Evaluate fitness $f(x_i^{new})$

Randonly choose a nest j away in

If $f(x_j^{new}) < f(x_j)$ then

Replace x_j with x_i^{new}

end if

end for

Replace x^{best} with x^{new}

Randomly choose a fraction ρ of worst nests and replace

with random feasible solutions

Rank nests, update best solution

End while

Return best nest as optimal gen settings

OP:-

Generator 1: 853.40 MW

Generator 2: 153.32 MW

Generator 3: 87.38 MW

Total Power: 1000.00 MW
Total Cost: \$ 4303.06

Return x^{best} as the best optimal solution

Code:

```
import numpy as np import
math
# Objective function to optimize (example: Sphere function)
def objective_function(x):      return np.sum(x**2)

# Lévy Flight distribution def
levy_flight(beta=1.5, size=1):
    sigma_u = (math.gamma(1 + beta) * np.sin(np.pi * beta / 2) /
math.gamma((1 + beta) / 2) * beta * (2 ** ((beta - 1) /
2)))**(1 / beta)      u = np.random.normal(0,
sigma_u, size)      v = np.random.normal(0,
1, size)      step = u / (np.abs(v) ** (1 /
beta))      return step

# Cuckoo Search Algorithm def cuckoo_search(objective_function, dim,
lower_bound, upper_bound, num_nests=25, max_iter=100, pa=0.25):
    # Initialize nests with random solutions within bounds      nests =
np.random.rand(num_nests, dim) * (upper_bound - lower_bound) + lower_bound
fitness = np.apply_along_axis(objective_function, 1, nests)
    # Initialize the best solution
best_nest_idx = np.argmin(fitness)      best_nest =
nests[best_nest_idx]      best_fitness =
fitness[best_nest_idx]

    # Iterate for a fixed number of generations or until convergence
for iteration in range(max_iter):      for i in range(num_nests):
        # Generate a new solution using Lévy flight
step = levy_flight(size=dim)      new_nest = nests[i] + 0.01
* step      new_nest = np.clip(new_nest, lower_bound,
upper_bound)
        # Evaluate the new solution
new_fitness = objective_function(new_nest)

        # If the new solution is better, replace the old solution
if new_fitness < fitness[i]:      nests[i] = new_nest
fitness[i] = new_fitness
```

```

        # Abandon the worst nests
for i in range(num_nests):
    if np.random.rand() < pa:  # Probability to abandon
nests[i] = np.random.rand(dim) * (upper_bound - lower_bound)
+ lower_bound
fitness[i] =
objective_function(nests[i])
    # Find the current best nest
best_nest_idx = np.argmin(fitness)
best_nest = nests[best_nest_idx]
best_fitness = fitness[best_nest_idx]

    # print(f"Iteration {iteration+1}, Best Fitness: {best_fitness}")
return best_nest,
best_fitness

# Example usage of Cuckoo Search

# Define the problem dimensions and bounds dim = 5
# Dimension of the solution space lower_bound = -5
# Lower bound of the search space upper_bound = 5
# Upper bound of the search space

# Run Cuckoo Search best_solution, best_fitness =
cuckoo_search(objective_function, dim, lower_bound, upper_bound,
num_nests=25, max_iter=100, pa=0.25)
print("Best Solution:
{best_solution}") print("Best Fitness:
{best_fitness}")

Output :
```

```

Best Solution: [0.64982748 0.55961241 2.01501756 0.93987275 0.31984962]
Best Fitness: 5.78140211553397
```

Program 5

Grey Wolf Optimizer (GWO)

Algorithm:

Date: 17/10/21	Page: 19
→ Grey Wolf Optimizer	
GWO scheduling	
Appn: Scheduling and Resource Allocation	
Wt. scheduling()	
Inital population of wolves (solution)	
π_i for $i=1 \dots n$	
- each π_i represents a task-to-machine assignment (vector of length N)	
- $\pi_i = \pi_i = [0, 1, 2, 1, 0, \dots]$ means task $\phi \rightarrow M_0$	
task $\phi \rightarrow M_1$, etc.	
Evaluate fitness of each π_i	
- for each solution compute makespan (maximise machine completion time)	
Identify Alpha (best), Beta (second best), Gamma (third best) wolves	
while ($t < No\text{ of Creations}$)	
for each wolf π_i in population:	
for each dimension j (task assignment):	
- Compute pos ⁿ update using Alpha, Beta, Gamma:	
$D_{alpha} = C_1 * \pi_i(\text{alpha}[j]) - \pi_i(\text{beta}) $	
$D_{beta} = C_2 * \pi_i(\text{beta}[j]) - \pi_i(\text{gamma}) $	
$D_{delta} = C_3 * \pi_i(\text{delta}[j]) - \pi_i(\text{gamma}) $	
$x_1^j = 2 * \pi_i(\text{alpha}[j]) - A_1 * D_{alpha}$	
$x_2^j = 2 * \pi_i(\text{beta}[j]) - A_2 * D_{beta}$	
$x_3^j = 2 * \pi_i(\text{delta}[j]) - A_3 * D_{delta}$	
$x_i^{new[j]} = \text{round}((x_1^j + x_2^j + x_3^j) / 3)$	
Update a, A, C	
Evaluate fitness of new positions	
Update alpha, Beta, Delta wolves	
$t = t + 1$	
Return alpha as best task scheduling	
OP:-	
Task Processing Times: [2 9 8 6 4 6 2 6 4 4]	
Best Machine Assignment: [0 0 2 1 2 1 1 2 0 1]	
Minimum Makespan Achieved: 18.0	

Code:

```
import numpy as np

# Objective function (example: Sphere
function) def objective_function(x):
return np.sum(x**2)

N, dim, T = 30, 10, 100 # Number of wolves, dimensions, iterations
lower_bound, upper_bound = -10, 10
wolves = np.random.uniform(lower_bound, upper_bound, (N, dim))

alpha_pos, beta_pos, delta_pos = np.zeros(dim), np.zeros(dim),
np.zeros(dim) alpha_score, beta_score, delta_score = float('inf'),
float('inf'), float('inf') for t in range(T): for i in range(N):
    fitness = objective_function(wolves[i]) # Evaluate fitness
if fitness < alpha_score:
    delta_score, delta_pos = beta_score, beta_pos.copy()
beta_score, beta_pos = alpha_score, alpha_pos.copy()
alpha_score, alpha_pos = fitness, wolves[i].copy() elif
fitness < beta_score:
    delta_score, delta_pos = beta_score, beta_pos.copy()
beta_score, beta_pos = fitness, wolves[i].copy() elif
fitness < delta_score:
    delta_score, delta_pos = fitness,
wolves[i].copy() a = 2 - t * (2 / T) for i in
range(N):
    r1, r2 = np.random.rand(dim), np.random.rand(dim)
A, C = 2 * a * r1 - a, 2 * r2 wolves[i] += A *
(abs(C * alpha_pos - wolves[i]) +
abs(C * beta_pos - wolves[i]) +
abs(C * delta_pos - wolves[i]))
    wolves[i] = np.clip(wolves[i], lower_bound,
upper_bound) print("Best Solution:", alpha_pos) print("Best
Score:", alpha_score)
```

Output :

```

Best Solution: [-1.28434275  1.94786008  0.82301541 -1.85113457 -2.08806377
3.74582237
0.84065243  0.8938704 -1.22271966 -0.29007149]
Best Score: 31.023829961456407

```

Program 6

Parallel Cellular Algorithms and Programs

Algorithm:

	<p style="text-align: right;">Date 07/11/25 Page 20</p> <p>→ Parallel Cellular Algorithms and Programs</p>
(i)	Define objective $f(x)$ to be minimized or maximized
(ii)	Set algorithm parameters: - grid size - num_iterations - neighborhood_type - alpha (diffusion coefficient) - mutation_rate
(iii)	Initialize each cell with a random position x_0 in the search space
(iv)	Evaluate $f(x_0)$ for each cell
(v)	for each iteration $t = 1$ to num_iterations: for each cell i in the grid: - Get neighbor cells based on neighborhood type - Find best neighbor soln x_{best} - Update cell's solution: $x_{i,new} = x_i + \alpha * (x_{best} - x_i) + \text{mutation rate} * \text{random noise}$
	Evaluate $f(x_{i,new})$ for all cells
	Replace x_i with $x_{i,new}$ if it improves fitness
	Track global best solution
(vi)	Output best solution found

Output :-

Iteration 0: Best value = 2.6008660
10: Best value = 0.089826
20: Best value = 0.061090
30: Best value = 0.061090
40: Best value = 0.024309
50: Best value = 0.024309
60: Best value = 0.019575
70: Best value = 0.009910
80: Best value = 0.009910
90: Best value = 0.009910

~~Best solution found:~~

$$\mathbf{x} = [0.0033020 \quad 0.00623017]$$

$$f(\mathbf{x}) = 0.00991042917967900$$

Code:

```
import numpy as np
random import
concurrent.futures
```

```

def rastrigin(x):      A = 10      return A * len(x) + sum([(xi ** 2 - A *
np.cos(2 * np.pi * xi)) for xi in x])

GRID_SIZE = (10, 10)
DIM = 2
RADIUS = 1
ITER = 100
BEST = None
def init_grid(size,
dim):
    return [[np.random.uniform(-5.12, 5.12, size=(dim,)) for _ in
range(size[1])] for _ in range(size[0])]
def
fitness(cell):
    return rastrigin(cell)
def update_state(grid, i, j,
radius):
    curr = grid[i][j]      fitness_curr = fitness(curr)      neighbors =
[grid[ni][nj] for dx in range(-radius, radius+1) for dy in range(-radius,
radius+1)                  if 0 <= (ni := i+dx) < len(grid) and 0 <= (nj
:= j+dy) < len(grid[0]) and (dx or dy)]      if neighbors:
        best_neigh = min(neighbors, key=fitness)
    return curr + 0.1 * (best_neigh - curr)      return
curr
def run_iteration(grid,
radius):
    new_grid = [[None for _ in range(len(grid[0]))] for _ in
range(len(grid))]      with
concurrent.futures.ThreadPoolExecutor() as ex:
    futures = [ex.submit(update_state, grid, i, j, radius) for i
in range(len(grid)) for j in range(len(grid[0]))]      for idx,
future in enumerate(futures):          i, j = divmod(idx,
len(grid[0]))          new_grid[i][j] = future.result()      return
new_grid
def
track_best(grid):
    global BEST
    best_cell, best_fitness = None, float('inf')
for row in grid:

```

```

        for cell in row:
            f = fitness(cell)                  if f <
best_fitness:                      best_fitness = f
best_cell = cell      if BEST is None or best_fitness
< fitness(BEST):
    BEST = best_cell
def parallel_cellular_algorithm():    global
BEST      grid = init_grid(GRID_SIZE, DIM)
for _ in range(ITER):           grid =
run_iteration(grid, RADIUS)
track_best(grid)          print(f"Best Fitness:
{fitness(BEST)}")
print("Best Solution:", BEST)
print("Best Fitness:", fitness(BEST))
parallel_cellular_algorithm()

```

Output :

```

Best Fitness: 2.4309484366586602
Best Fitness: 2.4309484366586602
Best Fitness: 0.0007801439196555293
Best Fitness: 0.0007801439196555293
Best Fitness: 0.0007801439196555293
Best Fitness: 0.0007801439196555293
Best Solution: [ 0.00129305 -0.00150346]
Best Fitness: 0.0007801439196555293

```

Program 7

Optimization via Gene Expression Algorithms

Algorithm:

→ Pseudo Code.

- Start
- Define fitness function
- Define parameters
- Create population
- Select mating pool
- Migrate on after mating
- Gene expression and evaluation
- Create
- Output best value

Output: (After 1000 generations)

Genes: [29.53, 29.82, 29.84, 28.57, 15.09,
21.83, 23.83, 30.81, 28.51, 26.27]

$x = 26.37$
 $f(x) = 695.45$

Generation limit reached

Code:

```

import numpy as np

# Define the mathematical function to optimize (example: minimize f(x) = x^2)
def optimization_function(x):
    return np.sum(x**2) # Modify this for other functions to optimize

# Parameters
POPULATION_SIZE = 50 # Number of individuals
GENE_LENGTH = 5 # Number of genes (dimensions of the problem)
MUTATION_RATE = 0.1 # Probability of mutation
CROSSOVER_RATE = 0.7 # Probability of crossover
GENERATIONS = 100 # Number of generations
SEARCH_SPACE = (-10, 10) # Range of values for genes

# Initialize Population
def initialize_population():
    return np.random.uniform(SEARCH_SPACE[0], SEARCH_SPACE[1],
                           (POPULATION_SIZE, GENE_LENGTH))

# Evaluate Fitness (lower is better for minimization)
def evaluate_fitness(population):
    fitness = np.array([optimization_function(ind) for ind in population])
    return fitness

# Selection (Roulette Wheel Selection)
def select_parents(population, fitness):
    # Convert fitness to probabilities (lower fitness is better)
    inverted_fitness = 1 / (fitness + 1e-6) # Avoid division by zero
    selection_prob = inverted_fitness / np.sum(inverted_fitness)
    selected_indices = np.random.choice(np.arange(POPULATION_SIZE),
                                        size=POPULATION_SIZE, p=selection_prob)
    return population[selected_indices]

# Crossover (Blend Crossover)
def crossover(parents):
    offspring = np.empty_like(parents)
    for i in range(0, POPULATION_SIZE, 2):
        p1, p2 = parents[i], parents[i+1] if np.random.rand() < CROSSOVER_RATE:
            alpha = np.random.rand() # Blending factor
            offspring[i] = alpha * p1 + (1 - alpha) * p2
            offspring[i+1] = alpha * p2 + (1 - alpha) * p1 else:
                offspring[i], offspring[i+1] = p1, p2
    return offspring

```

```

# Mutation (Random Perturbation) def
mutate(offspring):      for i in
range(POPULATION_SIZE):          if
np.random.rand() < MUTATION_RATE:
    mutation_point = np.random.randint(0, GENE_LENGTH)
offspring[i][mutation_point] += np.random.uniform(-1, 1)
    # Keep within search space
offspring[i][mutation_point] = np.clip(offspring[i][mutation_point],
SEARCH_SPACE[0], SEARCH_SPACE[1])      return offspring

# Gene Expression (Translate Genetic Code into Solutions) def
gene_expression(genes):
    # In this simple example, the genes directly represent the solution
return genes

# Main Algorithm def
gene_expression_algorithm():      #
Initialize population      population =
initialize_population()
best_solution = None      best_fitness =
float('inf')

    # Iterate through generations      for
generation in range(GENERATIONS):
        # Evaluate fitness      fitness =
evaluate_fitness(population)

        # Track the best solution
        current_best_idx = np.argmin(fitness)
if fitness[current_best_idx] < best_fitness:
best_fitness = fitness[current_best_idx]
best_solution = population[current_best_idx]

        print(f"Generation {generation+1}: Best Fitness = {best_fitness}")
        # Selection      parents =
select_parents(population, fitness)

        # Crossover
offspring = crossover(parents)

        # Mutation
offspring = mutate(offspring)

```

```

    # Gene Expression (not needed explicitly as genes represent
solutions)           population = gene_expression(offspring)
    print("\nOptimal Solution Found:")
print("Best Solution:", best_solution)
print("Best Fitness:", best_fitness)

# Run the algorithm if __name__ ==
"__main__":
    gene_expression_algorithm()

```

Output :

```

Generation 1: Best Fitness = 16.545885126119284
Generation 2: Best Fitness = 11.641082640808637
...
Generation 99: Best Fitness = 0.02233046748484963
Generation 100: Best Fitness = 0.02233046748484963

```

```

Optimal Solution Found:
Best Solution: [ 0.07226226 -0.11854791  0.03245473 -0.01236219  0.04299877]
Best Fitness: 0.02233046748484963

```