
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

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CERTIFICATE

This is to certify that the Lab work entitled “ Bio Inspired Systems (23CS5BSBIS)” carried out by **Sanchit Kashyap(1BM23CS298)**, 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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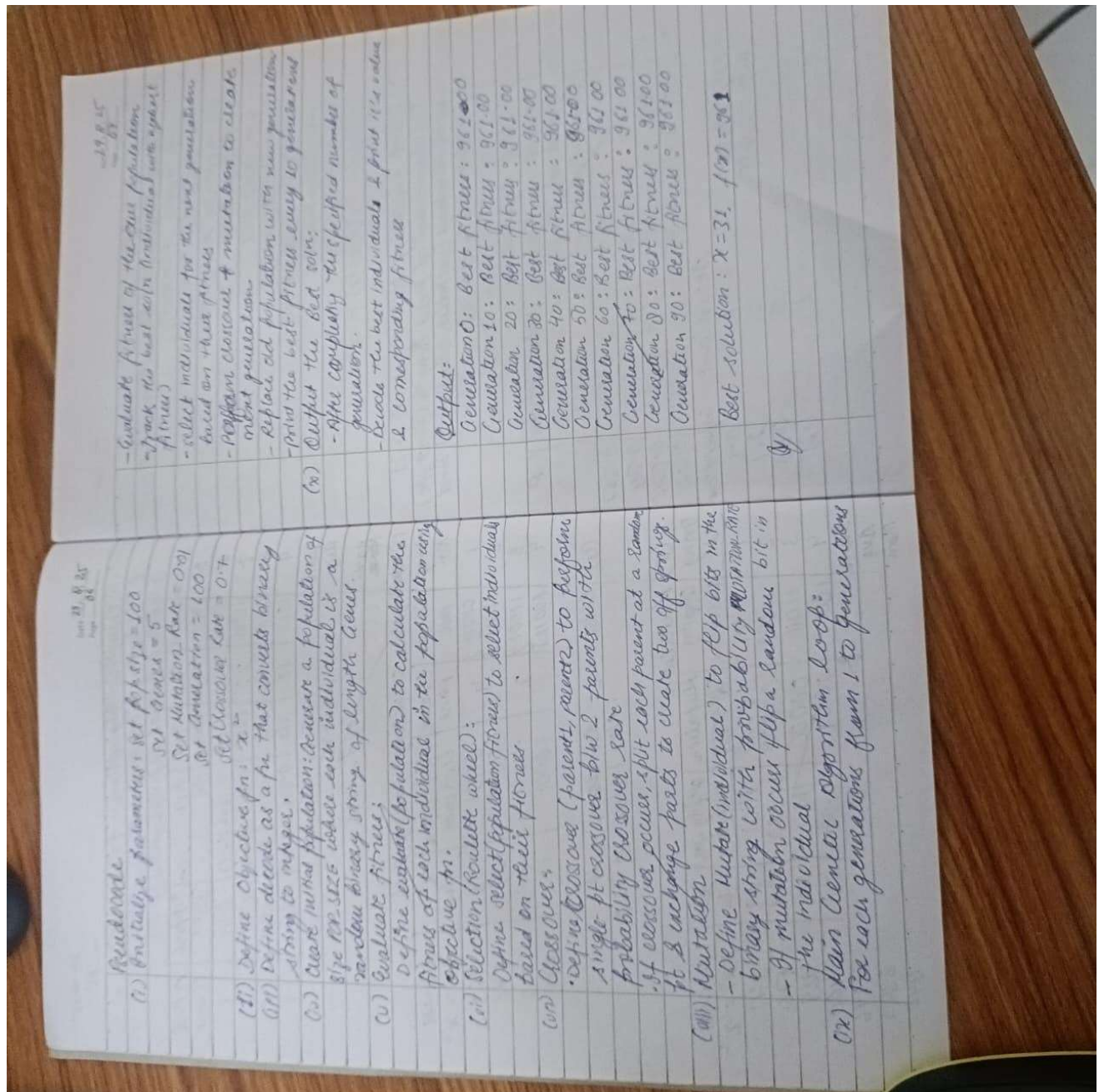
Github Link:

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:

15/09/25
Page 11

→ Particle Swarm Organization (PSO)

Pseudo code:-

- $P = \text{particle initialization()}$
for $i = 1 \rightarrow \text{max}$
for each particle p in P do
 $f_p = f(p)$
if f_p is better than $f(p_{\text{best}})$
 $p_{\text{best}} = p$
end
end for
 $g_{\text{best}} = \text{best } p \text{ in } P$
for each particle p in P do
$$U_i^{t+1} = U_i^t + C_1 U_i^t (p_{\text{best}}^t - p_i^t) + C_2 U_i^t (g_{\text{best}}^t - p_i^t)$$

inertia

personal
influence

social
influence

$$p_i^{t+1} = p_i^t + U_i^{t+1}$$

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	12
Page	12
Ant	1
City	2
Distance	2
Cost	0.5
Iteration	2
Iteration	0.125

Fitness
Value

2
2
0.5
2
125

Fitness
Value

125
125
125
125
125

→ Ant Colony Optimization for the Traveling Salesman Problem

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. proportional to $(\text{pheromone}[\text{current_city}][\text{city}]^\alpha) * (\text{heuristic}[\text{current_city}][\text{city}]^\beta)$

- Normalise 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).
- b. 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 $1 / \text{tour length}$:

$$\text{pheromone}[\text{city}_i][\text{city}_j] += Q / \text{tour length}$$
- c. Update best tour and best length if any ant found a better tour.
- d. Return best tour and best length

Output:-

cities = (0,0), (1,5), (5,2), (6,6), (8,3), (7,7)

~~n-ants = 20~~
n-ants = 20

n-iterations = 100

alpha = 1

beta = 5

rho = 0.5

initial-pheromone = 1

on the
red cities
blended tour

Iteration 1/10, Best length: 24.2828
2/10 : 24.2828

cities
ges:
distance

10/10, Best length: 24.2828

Best tour found: (1, np.int64(3), np.int64(5),
np.int64(2), np.int64(0))

all's tour

Best tour length: 24.282800682479206

length

$\frac{24}{10/10}$

if any

th

(7, 7)

Code:

```
import numpy as np
import 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] = 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

Pseudo Code:-

Cuckoo Search():

Initialize population of n host nests (solutions)

$q1 = 1, 2, \dots, n$

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

Find the current best solution x_{best} among the nests

While ($t < \text{MaxGenerations}$ OR stopping criterion)

For each nest $i = 1, 2, \dots, n$ do

Generate a new solution x_i^{new} by Levy flight from x_i

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

Randomly choose a nest j among n

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

Replace x_j with x_i^{new}

end if

end for

A fraction pa of worst nests are abandoned and replaced with new random solutions

Keep the best solution (x_{best}) among current popn.

Rank the nests and find the new x_{best}

$t = t + 1$

End while

Return x_{best} as the best optimal solution

APP: Optimal Power Flow

Initialize population of nests (generators settings, voltages, etc)

Evaluate power flow & cost for each nest

Find best nest

While (not stopping)

For each nest i

Generate new nest i new setting Levy flight

Run power flow, evaluate cost + penalty

If i new is better than a random nest j

Replace j with i new

End if

End for

Abandon a fraction pa of worst nests and replace with random feasible solutions

Rank nests, update best solution

End while

Return best nest as optimal gen settings

OP:-

Generators 1: 859.40 MW

Generators 2: 153.88 MW

Generators 3: 87.38 MW

Total Power: 500.00 MW

Total Cost: \$ 4303.06

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(f"Best Solution:
{best_solution}") print(f"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:

<p>→ Grey Wolf Optimizer</p> <p><u>Goal</u>: Scheduling and Resource Allocation</p> <p><u>App</u>: Scheduling and Resource Allocation</p> <p><u>GWO Scheduling()</u></p> <p>Initialize population of wolves (solutions)</p> <p>λ_i for $i=1, \dots, n$</p> <p>Each λ_i represents a task-to-machine assignment (vector of length n)</p> <p>Ex: $\lambda_i = [0, 1, 2, 1, 0]$ means task 0 \rightarrow M1</p> <p>task \rightarrow M1, etc.</p> <p>Evaluate fitness of each λ_i</p> <p>For each solution, compute makespan (maximum machine completion time)</p> <p>Identify Alpha (best), Beta (second best), Delta (third best) wolves</p> <p>While ($t < \text{MaxIterations}$)</p> <p>For each wolf λ_i in population:</p> <p>Compute post-update using Alpha, Beta, Delta:</p> $D_alpha = C1 * \lambda_alpha[i] - \lambda_i[i] $ $D_beta = C2 * \lambda_beta[i] - \lambda_i[i] $ $D_delta = C3 * \lambda_delta[i] - \lambda_i[i] $ $\lambda1 = \lambda_alpha[i] - A1 * D_alpha$ $\lambda2 = \lambda_beta[i] - A2 * D_beta$ $\lambda3 = \lambda_delta[i] - A3 * D_delta$ $\lambda_i_new[i] = \text{round}((\lambda1 + \lambda2 + \lambda3) / 3)$	<p>Update a, A, C</p> <p>Evaluate fitness of new positions</p> <p>Update Alpha, Beta, Delta wolves</p> <p>$t = t + 1$</p> <p>Return Alpha as best task scheduling</p> <p><u>OP:-</u></p> <p>Task Processing Times: [2 9 8 6 4 6 2 6 4 4]</p> <p>Best Machine Assignment: [0 0 2 1 2 1 1 2 0 1]</p> <p>Minimum Makespan Achieved: 180</p>
--	--

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)
```

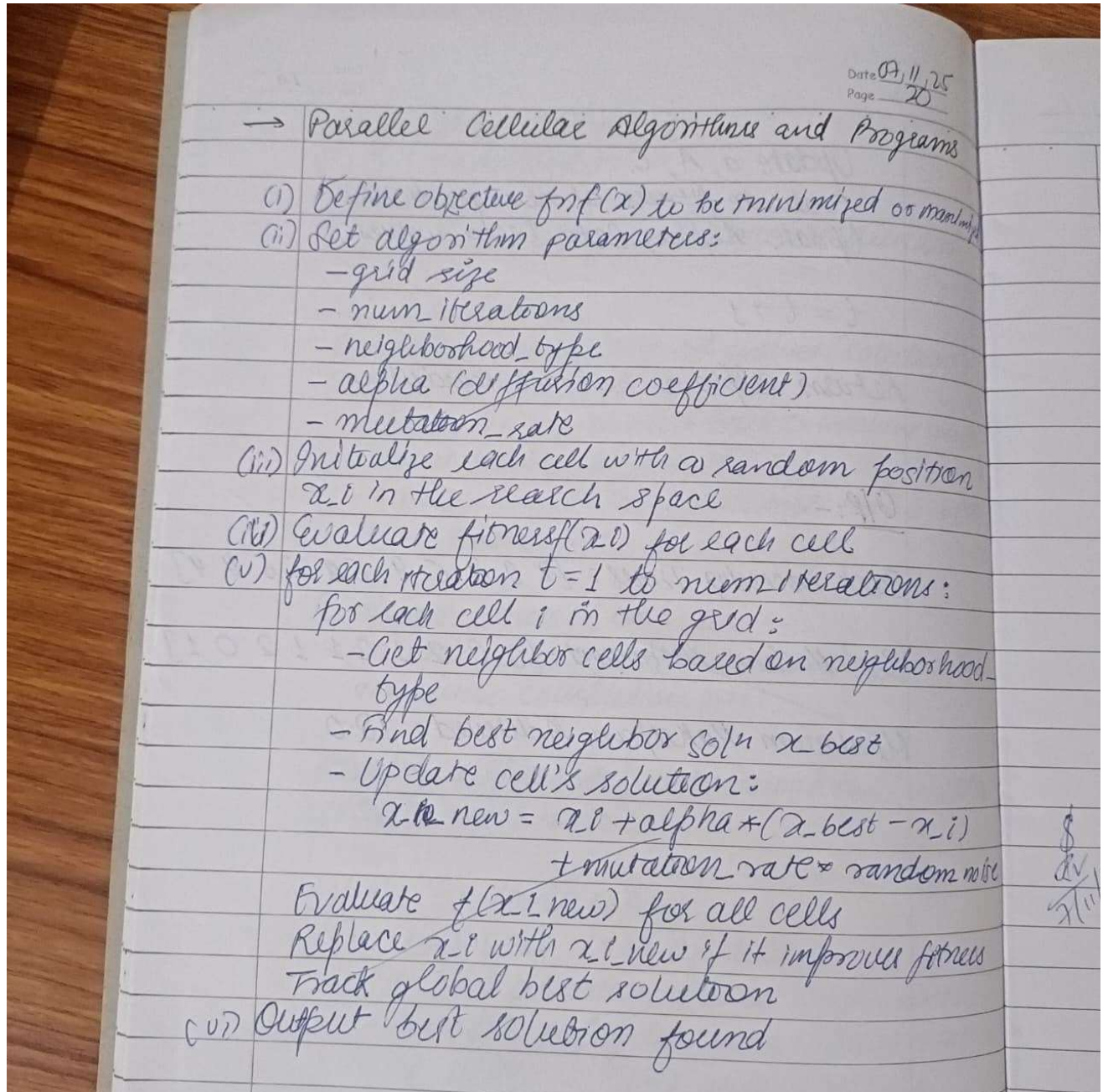
Output :

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:



Output:-

Iteration 0: Best value = 2.608660
 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:

$x = [0.003820 \quad 0.00623017]$

$f(x) = 0.00991042917967980$

\$
 25
 5/11/21

Code:

```

import numpy as np
import 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 Solution: [ 0.00129305 -0.00150346]
Best Fitness: 0.0007801439196555293

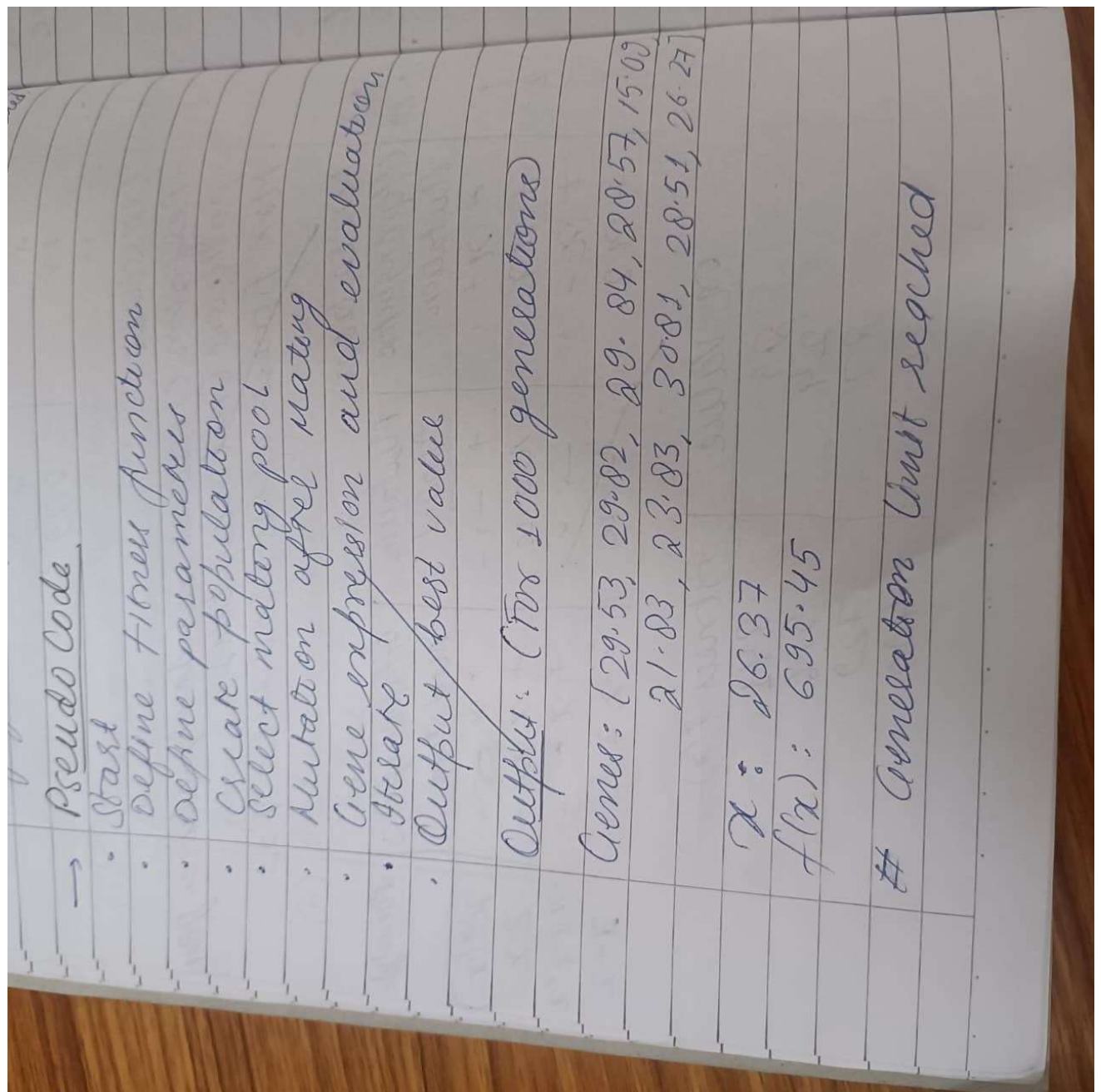
```

Program 7

Optimization via Gene Expression Algorithms

Algorithm:

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